

Roy F. Weston, Inc.
GSA Raritan Depot
Building 209 Annex (Bay F)
2890 Woodbridge Avenue
Edison, New Jersey 08837-3679
908-321-4200 • Fax 908-494-4021

5563

DATE:

March 24, 1995

TO:

Greg Powell, U.S. EPA/ERT Work Assignment Manager

THROUGH:

John N. Dougherty, REAC Geosciences Group Leader RMS for J. Doughe

FROM:

Ray Lewis, REAC Task Leader S. T Cor R.L

SUBJECT:

PIER DRUM, BRADENTON, FL

WORK ASSIGNMENT 0-098-TRIP REPORT

1.0 INTRODUCTION AND BACKGROUND

The U.S. Environmental Protection Agency (U.S.EPA) Region IV requested that the U.S. EPA/Environmental Response Team (ERT) conduct a geophysical survey to search for buried drums and collect and analyze soil, groundwater, surface water, and drum samples at the Pier Drum site in Bradenton, FL. The ERT activated the Response Engineering and Analytical Contract (REAC) on December 30, 1994, under Work Assignment 0-098, to provide personnel and equipment for a geophysical survey and drum, water, and soil sampling and analysis. Sam Getty, Michael Morganti, and Ray Lewis completed the sampling effort, and Stewart Sandberg and Jyotiranjan Kar performed the geophysical survey.

1.1 Site Description

The Pier Drum site is located in Bradenton Township, Manatee County, FL. The site is approximately 12 acres in size and is in a rural setting. Most of the property is open field, with a barn and sheds along the south side (Figure 1). In November 1994, the Florida Department of Environmental Protection (DEP) excavated 16 drums from two pits in the southeast corner of the site (Figure 2). The drums remain on-site near the excavation pits. The open pits subsequently filled with water, either from rain or groundwater or a combination of both.

1.2 Observations and Activities

Groundwater, surface water, sediment, and drum samples were collected during this sampling effort. In addition, geophysical monitoring and topographical surveying were conducted. Field work began on Monday, January 9, 1995 and concluded on Friday, January 13, 1995.

2.1 Sampling Activities

For all of the sampling activities, sample documentation was completed in accordance with ERT/REAC SOP # 2002, Sample Documentation, and ERT/REAC SOP # 4010, Chain of Custody (COC). The packaging and shipment of samples was completed in accordance with ERT/REAC SOP # 2004, Sample Packaging and Shipment. Field data sheets and COCs are provided in Appendix A.

2.1.1 Drum Sampling

Drums were sampled after they were opened with non-sparking beryllium tools. Samples were collected utilizing glass drum thieves and deposited into 32-oz glass jars for field screening. Drum samples were collected in accordance with ERT/REAC SOP # 2009, Drum Sampling. Drum samples were referenced according to the REAC's numbering system (REAC#1 through REAC#16) and the sample numbers correspond to Florida DEP numbers. Drums REAC#6, REAC#15, and REAC#16 could not be opened because they were filled to the top and the contents had solidified; therefore, they were not sampled. Drum REAC#12 was not sampled upon the instruction of the work assignment manager because it has the same contents as drum REAC#10.

2.1.2 Chemical Categorization Screening

Upon completion of the initial sampling, drum samples were chemically categorized using the Sensidyne HaztechTM Hazcat Kit (Hazcat) quick test methods. Quick test methods include the following individual tests: water reactivity/solubility, hairpin explosivity, oxidizer, acid/base pH, cyanide, chlorine hot wire, and organic peroxide. Samples which showed a positive response to the chlorine hot wire test were also screened with ChemetricsTM Quantichlor Chlorine in Waste Oil and DexsilTM Chlorin oil PCB test kits.

Each sample was tested for ignitability prior to the chemical testing with an ErdcoTM Setaflash Flash Point Tester (Setaflash). The Setaflash was set at 140 degrees Fahrenheit (°F) in order to characterize the sample as ignitable under Resource Conservation and Recovery Act (RCRA) regulations.

Samples were also characterized using a Spil-fyterTM Chemical test strip. Spil-fyter Chemical test strip consists of five colorimetric tests: an acid/base risk test; an oxidizer risk test; a fluoride risk test; an organic solvent/petroleum distillate risk test; and an iodine, bromine, and chloride risk test.

Upon completion of chemical categorization screening, six composite samples were sent out for further laboratory analyses. Composite samples 00911,00915, and 00916 were sent for TCLP VOA, TCLP BNA, TCLP PCB/Pesticides, and TCLP Metals analyses. Samples 00912, 00913, and 00914 were split and sent to two different laboratories for TCL VOC, BNA, and PEST/PCB, TAL metals, percent moisture, percent sulfur, percent chlorine, British Thermal Units (BTU), and ash content analyses.

The following is a description of the field tests that were used to characterize the samples.

Water Reactive/Solubility Test:

A small amount of sample was added to 1/2 inch of water in a test tube. If the sample did not effervesce, a stopper was placed in the test tube. The test tube was then shaken vigorously. After allowing time for a reaction to occur, the results were recorded. This test determined whether a sample is water reactive or soluble in water.

Oxidizer Test:

Two or three drops of acid test solution were placed on the oxidizer test paper and the sample was touched with the paper. An oxidizer turns the paper blue to black. If the test was positive, the pH of the sample was tested. If the pH was less than or equal to 2, the sample was classified as an oxidizer and no other tests needed to be performed. If the pH was greater than 2, the sample was tested for ignitability.

If the sample was ignitable, then it was classified as an organic peroxide. If it failed the ignitability test, the sample was classified as an oxidizer.

If the oxidizer test proved negative, the pH of the sample was tested. Samples with a pH greater than or equal to 12 were tested for CN. If the pH was between 2.5 and 12, the sample was tested for ignitability. For a pH less than 2.5, the sample was classified as an acid and the test ended.

Ignitability Tests:

The following tests were performed to determine the ignitability of the sample:

• Hairpin Test:

The hairpin test is a test for explosiveness. A grain-size solid or a drop of liquid sample was placed on a watch glass. A hairpin was heated until red hot, and then touched to the sample. If the sample ignited into flames, the sample was classified as an explosive.

If there was no reaction from the hairpin test, a lit match was used to ignite the liquid in a watch glass.

Observations: Extremely Flammable: Flame jumps from match to

liquid.

Flammable:

Stays ignited after

as a

removing match.

Combustible:

)----i--- ------

Requires match

wick.

Non-Flammable:

Will not ignite.

· Setaflash Test:

The Setaflash tester was set at 140 °F in order to characterize the liquid as ignitable under RCRA regulations. A small amount of each sample was placed into the chamber of the instrument and heated to 140°F. Once the sample amount reached 140°F, a flame was introduced to the headspace of the chamber. If a flash occurred, the sample was characterized as ignitable. The results were recorded in the log book as "flash" or "no-flash," with flash indicating that the sample was ignitable.

• Cyanide Test:

A small amount of CN test 2 solution was added to a test tube containing 1/4 inch of CN test 1 solution. Approximately 1/2 inch of the sample was added to another test tube. The CN test solution was added to the sample solution and then an acid test solution was added. A deep prussian blue color indicated the sample contained cyanide. A positive result for CN ended the test and the sample was classified as a CN. A negative result for CN ended the test and then the sample was classified as caustic.

• Chlorine Hot Wire Test:

A chlorine hot wire was heated in a torch flame until there was no green flame. The wire was then allowed to cool. The wire was placed into a test tube containing the sample and reheated in a torch flame. A green flame indicated the sample contained chlorine, amine, nitrate, or ammonium salt.

• Organic Halogenated Solvent Test:

Based on the results of the combustibility test, the Setaflash test, the chlorine hot wire test and the water reactive/solubility test, an assumption was made as to whether or not the sample was an organic halogenated solvent.

Clor-N-Oil test:

The Clor-N-Oil test was used to check samples for PCBs. A precise amount of sample was placed into a tube. A colorless ampule containing a catalyst was broken and thoroughly mixed with the sample. A second ampule containing metallic sodium was broken and the sodium, activated by the catalyst, stripped chlorine from the sample forming sodium chloride. A buffer solution in water was added to the sample which neutralized the excess sodium and extracted the sodium chloride into the water. The water layer was then separated from the sample.

An ampule containing a precise amount of reagent was broken and mixed with the water. An indicator ampule was then broken and mixed. The color of the mixture depended on the amount of PCBs (chlorine) in the sample.

The Clor-N-Oil test kit works on the principle of chloride determination. Since PCBs contain chlorine, the test kit is able to detect them. However, the test cannot distinguish between any other chlorine containing compound such as trichlorobenzene which may also be in transformer oil. This may cause a result which is known as a "false positive"; i.e., the oil may indicate the presence of over 50 ppm PCBs, but when analyzed by a gas chromatography, it will show somewhat less than 50 ppm.

Quanti-Chlor •1

In addition to and in support of a positive result with a Clor-N-Oil test, a Quanti-Chlor \bullet test kit was used. The Quanti-Chlor test yields quantitative results in micrograms per gram ($\mu g/g$) total chlorine in waste oil. Organic chlorine is converted to chloride with naphthalene, dig-lyme, and metallic sodium. The chloride was extracted into an aqueous buffer and titrated with mercuric nitrate. Diphenylcarbazone was the endpoint indicator. Titrets \bullet^2 , hand-held titration cells, were used in the mercuric nitrate titration.

Spil-fyter Chemical Classifier Test Strip:

In addition to the field screening tests, a Spil-fyter chemical classifier test strip was used on each sample. The results of the Spil-fyter strips were used as a secondary test to help support the data from the HAZCAT tests. Some interferences that the Spil-fyters have are concentrated acid solutions that tend to destroy the indicators, heavy oils and opaque solutions mask the colors, and light weight organic solvents cause color bleeding. The detection limits of the Spil-fyter are:

- 1 milligram/liter (mg/1), Chlorine,
- 3 mg/1, Hydrogen Peroxide,
- 20 mg/L, Fluoride and
- 10 mg/L, Gasoline.

2.1.3 Groundwater Sampling

Four temporary monitor wells were installed using hand augers (Figure 1). The PVC screens were installed approximately 2-3 ft below the water table. Prior to sampling, each monitor well was developed by purging until clear water (< 50 Nephelometric Turbidity Units) was produced. Well development was performed in accordance with U.S. EPA/ERT/REAC SOP # 2044, Well Development. Peristaltic pumps were used to develop and sample the wells. Between 20 and 25 gallons of water were purged from each well during development. Periodically during development, pH, dissolved oxygen, temperature, and conductivity were monitored using the Omega Water AnalyzerTM. Turbidity was monitored using the Lamont Turbidity MeterTM. The instruments were calibrated daily to ensure accurate measurements. Temperature, pH, dissolved oxygen, conductivity, and turbidity, were measured in each well (Table 1).

Depth to water measurements were taken after development of each well. Water level measurements were recorded in accordance with ERT/REAC SOP # 2043, Water Level Measurement. These measurements were used to determine groundwater elevation (Table 2) and flow direction.

¹ Quanti-Chlor • is a registered trademark of CHEMetrics, Inc.

² Titrets • is a registered trademark of CHEMetrics, Inc.

Groundwater samples were collected in accordance with ERT/REAC SOP # 2007, Groundwater Well Sampling. The groundwater samples were analyzed for Target Compound List (TCL) volatile organic compounds (VOCs), base, neutral, and acid extractable compounds (BNAs), pesticides/polychlorinated biphenyls (PEST/PCBs), and Target Analyte List (TAL) metals. Both filtered and unfiltered samples were analyzed for TAL metals.

2.1.4 Surface Water Sampling

Two surface water samples were collected from the excavation pits in accordance with ERT/REAC SOP # 2013, Surface Water Sampling. The surface water samples were analyzed for TCL VOAs, BNAs, PEST/PCBs, and TAL metals.

2.1.5 Sediment Sampling

Three sediment samples were collected from the excavation pits using hand augers. The sediment samples were collected in accordance with ERT/REAC SOP # 2016, Sediment Sampling. These samples were analyzed for TCL VOCs, BNAs, PEST/PCBs and TAL metals. In addition, these samples were analyzed for Toxic Characteristic Leaching Procedure (TCLP) for the same parameters.

2.1.6 Geophysical Survey

A survey was performed to detect and delineate geophysically anomalous areas indicative of buried drums within a 12-acre portion of the site. Variations of the earth's local magnetic field intensity occur in the vicinity of ferromagnetic objects. It is common practice to measure this field, alone or in combination with its vertical gradient, to locate buried ferromagnetic objects (such as steel drums). Empirical studies have shown that an individual 55-gallon steel drum can be detected when buried in relatively nonmagnetic material to depths of up to 10 feet.

Electromagnetic instruments can be used to detect anomalous conductivity enhancements such as those due to buried metallic objects (such as steel drums). Scattering of the quadrature-phase component of the induced electromagnetic field is used to determine the bulk electrical conductivity, or terrain conductivity, of the shallow subsurface. Enhancements of the inphase component of the scattered electromagnetic field result from the presence of relatively good conductors, such as buried metallic objects.

Geophysical methods used in this investigation included electromagnetics and magnetics. A GEM Systems model GSM-19 magnetometer/gradiometer was used to obtain the earth's total magnetic field intensity and vertical magnetic field gradient at geophysical grid stations. In addition, a Geonics EM-31 terrain conductivity instrument was used to obtain terrain conductivity and inphase readings at these same locations. A brief description of these methods follows.

A geophysical grid was established using a compass and a measuring tape, consisting of traverses spaced 20 feet apart, upon which stations were located at 5-foot intervals. The geophysical grid is shown on data plots in which station locations are posted with a "+"symbol (Figure 3). Line orientations were magnetic north-south, and wooden stakes were used to indicate the locations of the 0 North grid position on each line.

Data obtained each day were downloaded into a portable computer where profiles were viewed for preliminary data inspection and interpretation. Hard copy plots of specific traverses were produced to aid in interpretation. A preliminary interpretation was available the following morning.

Electromagnetics (EM-31): The electromagnetic method (EM) is a geophysical technique based on the physical principle of inducing and detecting electrical current flow within the earth, including any conductors contained (buried) therein. The EM-31 is a fixed-coil (approximately 10-foot separation) electromagnetic instrument that generates an electromagnetic field at a known frequency. The primary coil transmits the electromagnetic field and the receiving coil measures the change in electric and magnetic components (secondary electromagnetic field) of the field after propagation through the earth. The EM-31 measures two quantities: the inphase and quadrature phase components of the secondary electromagnetic field. The inphase measurement is more sensitive to good conductors (metallic objects) than the quadrature phase component. The quadrature phase measurement is commonly used to measure directly the bulk conductivity (or its inverse, resistivity) of lithologic materials. For this reason, the inphase component was closely reviewed for anomalies indicative of buried drums. Also, the quadrature phase component was useful for delineating changes in the conductivity due to excavation and subsequent burial.

Prior to use of the EM-31, the instrument was calibrated according to the manufacturer's recommendations. A data logger was used to record both inphase and quadrature phase measurements as well as all pertinent site features (e.g., metallic debris, power lines, buildings, etc.) to ensure quality data and assist in anomaly delineation.

The EM data was obtained at 5-foot intervals along traverses spaced 10 to 40 feet apart, depending upon specific site characteristics. A series of these traverses formed a data collection grid. The end points of each traverse were marked for subsequent anomaly location using stakes, flagging, or spray paint.

Anomalous responses resulting from metallic objects located on the surface were plotted in profile and contour form. Profile plots were used for interpretation on a per line basis, and contour plots were used to visualize spacial variations in instrument response over a large area.

<u>Magnetics</u>: The objective of the magnetic survey was to map magnetic field anomalies and thereby locate their ferromagnetic sources, in this case steel drums. The lines and sample nodes were spaced in the identical locations as those of the EM transects. Both magnetic and magnetic gradient data were electronically recorded.

The magnetic method involves precisely measuring the earth's magnetic field. A magnetometer is used to measure variations in this field over an area of interest. Local variations, or anomalies, in the earth's magnetic field are caused most often by nearby concentrations of ferromagnetic material. In this specific instance, a metallic cylinder, cement pad, steel fences and gates, as well as other surficial metallic debris, were the cause of the detected anomalies. The magnetometer used was the GEM Systems, GSM-19B Overhauser Effect Gradiometer.

2 2 3037

A magnetic base station was established in an area of low magnetic gradient (approximately ± 5 nanoTesla/meter) and measurements were periodically taken at this location during the magnetic survey to monitor the diurnal variation of the earth's magnetic field. As above, all pertinent site features (e.g., metallic debris, metal trailer, metal tank, steel gate, cement pad, steel fence, etc.) were recorded in a field notebook to ensure quality data and assist in anomaly delineation. All magnetic and gradient data were electronically recorded in the GSM-19B control console/data logger.

Total field magnetic data normally include the effects of diurnal variations of the earth's magnetic field. The diurnal variations were closely monitored by repeated observation of the total magnetic field every half hour.

2.1.7 Topographical Surveying

A topographic survey was conducted using a TopconTM electronic total station. The total station uses infrared light to calculate distances and angles between points. The survey was used to determine the relative locations and elevations of the geophysical grid, sampling locations, and other site features such as buildings, fences, and roads. This data was used to produce the geophysical data and the maps included in this report and, when combined with the water level measurements from the wells, to determine the direction of groundwater flow.

3.0 RESULTS

3.1 Sampling Activities

3.1.1 Chemical Categorization Screening

Drum contents were categorized using several field screening methods (Table 3). Based on the results of the screening, samples were composited into three solid samples and three liquid samples. The solid samples were analyzed for TCL/TAL and TCLP analytes. The liquid composites were analyzed for disposal characteristics and TCL analytes (Table 4).

Appendix A contains the field notes and Appendix B contains the results of additional field screening methods.

3.1.2 Drum Samples

The results of the analysis of the three liquid composites indicate high levels of VOCs, especially actione and styrene (Table 5). No BNAs were detected in any of these liquid samples. All three drum samples contained elevated levels of sodium, potassium, manganese, magnesium, and cobalt. The results of the disposal characterization analyses indicate that samples REAC #2 and REAC #9 have high BTU levels and low percent moisture where as sample REAC #4 has a low BTU and high percent moisture (Table 6). Percent ash, chloride, and sulfur were relatively low and consistent in all three samples. REAC #2 and REAC #4 were overpacked due to poor condition.

2 2 16:3

The results of the TCLP analysis for the drum solids also indicate that high concentrations of VOCs are present. In addition, two BNAs (benzylalcohol and benzoic acid) and two metals (sodium and calcium) were detected. Concentration levels of PEST/PCBs were very low level or non-detect (Table 7).

3.1.3 Groundwater Samples

Elevated concentrations of metals and VOCs were detected in groundwater samples. However, no concentrations were detected above the Maximum Concentration Limit (MCL) for those analytes with an established MCL. All samples are below federal and Florida state MCLs. There was very little difference in the concentration levels of the filtered and unfiltered samples analyzed for metals. This indicates that a majority of the metals in the groundwater are dissolved. PEST/PCBs were not detected in groundwater samples (Table 8).

3.1.4 Surface Water Sampling

PEST/PCBs were not detected in any surface water sample. In addition, one BNA (di-n-butylphthalate) and one VOC (acetone) were at low concentrations. The metals analyses results indicate that the majority of the metals in the surface water samples are dissolved in the water because there was little difference between the filtered and unfiltered sample concentrations (Table 9). Sodium, magnesium, calcium, and aluminum were detected at elevated levels. No sample concentrations were detected above MCLs.

3.1.5 Sediment Samples

Phthalates were the only BNAs detected in the sediment samples. However, these compounds were also detected in the blank. Metals and low levels of VOCs were also detected in sediments (Table 10).

The results of the TCLP analyses also indicate phthalates and benzylalchohol were detected. Of the BNAs detected, di-n-butylphthalate, butylbenzylphthalate, and bis(2-ethylhexyl)phthalate were also detected in the blank. Some metals were detected in low concentrations. No PEST/PCBs or VOCs were detected in any of the samples (Table 11).

3.2 Geophysical Survey

3.2.1 Geophysical Results

Contour plots of magnetic field gradient, total magnetic field, EM-31 inphase, and EM-31 terrain conductivity are depicted in Figures 4 through 7, respectively. Note that some contours and contour line extensions were plotted where no data points are located, and are therefore inappropriate for interpretation. All contour lines drawn greater than 20-feet from a data location (the "+"symbols) are considered interpolations.

Electromagnetic site features, observed surficial metallic objects (such as a trailer), a metal tank, debris pile consisting of rubber tires and sheet metal, and observed drums are evident in the EM-31 inphase contour plot (Figure 6). All major metallic-type anomalies can be accounted for by surficial metal. A partially buried drum was observed within a shallow pit on Line 60E, Station 150N, which is the location of all of the excavated drums.

2 2 0000

A metal automobile gas tank was noted on Line 100W, Station 125N and is apparent in the data as the peak in the EM-31 inphase (see Appendix C). The peak in the EM-31 inphase data, as shown in EM Line 60E, Station 150N (Appendix C) corresponds to the interpreted cache of drums. The magnetic gradient field corresponding to Line 60E, Station 150N indicates a low peak anomaly. The pattern depicting the surface debris pile on line 60E, Station 50N (Appendix C) indicates a high peak for the EM-31 inphase component and a low trough for the magnetic gradient component.

The diurnal variations may be neglected after careful observations of the total magnetic field at the base station and comparison with a strong anomaly of about 47500nT due to drums exposed in the pit on Line 60E, Station 150N (Figure 5). Therefore, the diurnal variation amplitude is relatively insignificant, indicating that diurnal corrections to these data are not necessary for the purposes of identifying large metallic responses.

3.3 Groundwater Elevation and Flow

The depth to water measurements were recorded after the development of each well. Based on the elevations of the groundwater, the flow direction was determined (Figure 8). Groundwater flow appears to be flowing in an easterly direction.

4.0 Conclusions

Sixteen drums were sampled to determine disposal characteristics. Based on the results of the field tests, the drum samples were composited into three liquid and three solid samples. Results of the analyses indicate elevated levels of volatile organics and metals.

Surface water, groundwater, and sediment samples were also collected. No analyte exceeded established MCLs for the groundwater and surface water samples. Low levels of VOCs, metals, and BNAs were detected in sediment samples.

Based on the results of a geophysical survey, all anomalies can be accounted for based on surface metal (known as drums and structures). No other anomalies were noted which would indicate additional buried drums.

Tables

TABLE 1 FINAL GROUNDWATER FIELD PARAMETERS PIER DRUM SITE BRADENTON, FLORIDA JANUARY 1995

	Turbidity	(NTU)
	Conductivity	(m2/cm)
Parameters	Dissolved	Oxygen (ppm)
	Temperature	(2)
		Hd
	e Installed Date developed	and Sampled pH
	Date Installed	
	Well ID	

NTU = Nephelometric Turbidity Units μS = micro Siemens (10 -6 Siemens)

2

2 1 9002

TABLE 2 GROUNDWATER ELEVATION DATA PIER DRUM SITE BRADENTON, FLORIDA JANUARY 1995

	SURFACE	DEPTH TO	GROUNDWATER
	ELEVATION	WATER	ELEVATION
LOCATION	(ft)	(ft)	(ft)

MW-N	99.71	1.88	97.83
MW-S	100.45	2.58	97.87
MW-E	99.63	1.91	97.72
MW-W	100.75	2.81	97.94

All elevations are relative to the site.

TABLE 3 CHEMICAL CATEGORIZATION SCREENING PIER DRUM SITE BRADENTON, FLORIDA 10 JANUARY 1895

7) [T	Ι	Ι	Γ			i	1		Ţ	_			<u> </u>	<u> </u>
ΨN	4,5,6	N/A	A/N	A/N	N/A	A/A	A/A	A/A	Ψ/Z	A/A	A/A	Y/X	Y/X	A/X	A/N
N/A	840-2400 ppm	A/N	A/N	K/N	N/A	N/A	A/A	A/A	A/N	A/N	A/X	A/N	A/A	A/A	A/A
z	+	A/X	A/A	A/A	A/A	A/N	A/X	A/X	N/A	N/A	N/A	N/A	N/A	N/A	W/A
N/A	S&F	S	N/A	L.	A/A	A/A	S	N/A	S	S	A/A	N/A	ц.	N/A	A/A
N-C	O-Z	N-C	N-C	N-C	N/A	N-C	N-C	Cloudy White	O-N	N-C	N/A	N-C	N-C	N/A	A/X
d	۵	z	1	Z	A/A	z	Q	1	Z	Z	N/A	z	Z	N/A	A/A
5	7	7	7	A/A	Y/X	7	7	7	7	7	A/A	7	7	A/A	V/V
FLASH	NO-FLASH	NO-FLASH	NO-FLASH	NO-FLASH	A/A	NO-FLASH	NO-FLASH	NO-FLASH	NO-FLASH	FLASH	¥/X	NO-FLASH	NO-FLASH	¥/X	A/A
×	:	×		×	∢ Z	×	×	×	×	×	∀/Z	×	×	₹ Z	Α/X
	×		×		∀\Z			×			ĕ, Z			ĕ,Z	¥ X
17E	17E	17E	17.6	17E	17E	17E	17E	17E	17E	17E	17E	17E	17E	17E	17E
01	02	03	40	0.5	90	07	80	60	0	=	12	13	4	15	16
REAC#1	REAC#2	REAC#3	REAC#4	REAC#5	REAC#6	REAC#7	F! AC#8	HEAC#9	REAC#10	HEAC#11	REAC#12	REAC#13	REAC#14	REAC#15	REAC#16
501 I	502	503	504	505	909	/*:	6	ن	019		512	513	_	615	616
	REAC#1 01 17E X FLASH 5 P N-C N/A N/A	REAC#1 01 17E	REAC#1 01 17E	REAC#1 01 17E	REAC#1 01 17E X FLASH 5 P N-C N/A N/A N/A REAC#2 02 17E X NO-FLASH 7 P N-C S&F + 840-2400 ppm REAC#3 03 17E X NO-FLASH 7 N N-C S N/A N/A N/A REAC#4 04 17E X NO-FLASH 7 T N-C N/A N/A	REAC#1 01 17E X FLASH 5 P N-C N/A N/A N/A REAC#2 02 17E X NO-FLASH 7 P N-C S&F + 840-2400ppm REAC#2 03 17E X NO-FLASH 7 N N-C S N/A N/A REAC#2 04 17E X NO-FLASH 7 N N N/A N/A N/A REAC#6 05 17E X N/A N/A N/A N/A N/A REAC#6 06 17E N/A N/A N/A N/A N/A	REAC#1 01 17E	Neac#1 01 17E	Neac#1 01 17E	REAC#1 01 17E X FLASH 5 P N-C N/A N/A N/A REAC#2 02 17E X NO-FLASH 7 P N-C S&F + 840-2400ppm REAC#2 03 17E X NO-FLASH 7 T N-C S N/A N/A N/A REAC#4 04 17E X NO-FLASH 7 T N-C F N/A N/A N/A REAC#5 05 17E N/A N/A N/A N/A N/A N/A N/A REAC#6 06 17E N/A N/A N/A N/A N/A N/A N/A REAC#6 06 17E X NO-FLASH 7 N N-C S N/A N/A REAC#6 08 17E X NO-FLASH 7 P N-C S N/A N/A REAC#9	REAC#1 01 17E X FLASH 5 P N-C N/A N N/A REAC#2 02 17E X NO-FLASH 7 P N-C S&F + 840-2400ppm REAC#2 02 17E X NO-FLASH 7 T N-C S N/A N/A REAC#2 04 17E X NO-FLASH 7 T N-C F N/A N/A REAC#5 05 17E N/A N/A N/A N/A N/A N/A REAC#6 06 17E N/A N/A N/A N/A N/A REAC#6 07 17E X NO-FLASH 7 N N-C S N/A N/A REAC#9 08 17E X NO-FLASH 7 N N-C S N/A N/A REAC#9 09 17E X NO-FLASH 7	No	REAC#1 01 17E X FLASH 5 P N-C N/A N/A N/A REAC#2 02 17E X NO-FLASH 7 P N-C S&F + 840-2400ppm REAC#2 03 17E X NO-FLASH 7 N N-C S N/A N/A REAC#4 04 17E X NO-FLASH 7 N N-C F N/A N/A REAC#6 06 17E N/A N/A N/A N/A N/A N/A REAC#6 06 17E N/A N/A N/A N/A N/A N/A REAC#6 06 17E X NO-FLASH 7 N N/A N/A N/A REAC#9 09 17E X NO-FLASH 7 N N N N/A N/A REAC#10 10 17E X NO-FLASH 7	Neac#1 01 17E	NO-FLASH 01 17E

TP = Department of Environmental Protection

1/A = Non Applicable
1 - C = No Change
2 = Sinks, F = Floats
4 = Organic Solvent, Petroleum, Distlate Risk; 5 = Iodine, Bromine, Chlorine Risk
T = Total Solubility, P = Parcial Solubility; N = No Solubility
NOTE: Drums #6, #15, #16 could not be opened, and drum #12 was determined to have the same contents as drum #10.

2 1 3004

TABLE 4 DRUM SAMPLE COMPOSITE SCHEME PIER DRUM SITE BRADENTON, FLORIDA JANUARY, 1995

Composing Scheme										
Composite		Sample	Analysis							
ID	Sample ID	Phase								

00912	REAC #9*	liquid	Disposal & TCL/TAL
00913	REAC #2	liquid	Disposal & TCL/TAL
00914	REAC #4	liquid	Disposal & TCL/TAL
00911	REAC #1,8,9*,11	solid	TCL (TCLP)
00915	REAC #5	solid	TCL (TCLP)
00916	REAC #3,7,10,13,14	solid	TCL (TCLP)

^{*} Drum #9 had both a liquid and a solid layer
Disposal Parameters include: % moisture, % sulfur, % chloride, BTU, Ash content
TCL - refers to the Target Compound List for VOC, BNA, and Pest/PCB
TAL - refers to the Target Analyte List for metals

0025

TABLE 5 SUMMARY OF RESULTS FOR DRUM LIQUID COMPOSITE SAMPLES PIER DRUM SITE **BRADENTON, FLORIDA JANUARY 1995**

SAMPLE ID	00912		0091	13	00914			
SAMPLE LOCATION	REAC	#9	REAC	#2	REA	C#4		
	Conc	MDL	Conc	MDL	Conc	MDL		
PARAMETER	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L		
METALS								
Aluminum	690	100	ND	100	25000	100		
Arsenic	ND	10	ND	10	22	10		
Barium	29	10	23	10	240	10		
Cadmium	7.0	4.0	34	4.0	ND	4.0		
Calcium	410*	0.20*	400*	0.20*	37*	0.20*		
Chromium	80	10	120	10	47	10		
Cobalt	7300	10	40000	10	900	10		
Copper	520	10	150	10	660	10		
Iron	18*	0.05*	360*	0.05*	310*	0.05*		
Lead	61	10	1200	10	38	10		
Magnesium	8000	1000	8000	1000	10000	1000		
Manganese	3000	4.0	5700	4.0	1500	4.0		
Nickel	66	20	850	20	30	20		
Potassium	12000	4000	35000	4000	12000	4000		
Sodium	21000	1000	71000	1000	29000	1000		
Vanadium	ND	10	ND	10	32	10		
Zinc	1900	10	3600	10	330	10		
PEST/PCB's								
a-BHC	0.16	0.11	ND	0.11	ND	0.11		
Aldrin	ND	0.11	0.18	0.11	ND	0.11		
VOC's								
Acetone	130000*	10*	230000*	10*	14000*	20*		
2-Butanone	ND	20000	ND	20000	89000	4000		
Benzene	11000	5000	12000	5000	ND	1000		
Toluene	5400	5000	4100J	5000	9600	1000		
Ethylbenzene	21000	5000	88000	5000	37000	1000		
p&m Xylene	16000	5000	51000	5000	ND	1000		
o-Xylene	8400	5000	ND	5000	800J	1000		
Styrene	1200000	5000	16000000	5000	26000	1000		
Isopropylbenzene	6500	5000	21000	5000	ND	1000		
n-Propylbenzene	7600	5000	17000	5000	ND	1000		
4 O 4 Tales Alexable a second	17000	5000	22222			1		

5000

5000

23000

ND

45000

7200

1000

1000

MDL = Method Detection Limit

1,2,4 - Trimethylbenzene

ND = Not Detected

Naphthalene

* = Concentration and MDL reported in milligram per liter (mg/L)

NOTE: Mercury analysis was not performed due to insufficient sample size.

5000

5000

17000

3400J

2 2 0006

TABLE 6 RESULTS OF ANALYSIS FOR DISPOSAL CHARACTERISTICS ON DRUM LIQUIDS PIER DRUM SITE BRADENTON, FLORIDA JANUARY 1995

	Composite ID REAC # 9 REAC # 2 REAC # 4									
	REAC # 9	REAC # 2	REAC # 4							
Parameter	ID # 00912	ID # 00913	ID # 00914							

Heat Value (BTU/lb)	3379	5470	+50
Percent Ash Content	0.2	0.26	+0.1
Percent Chloride	0.33	0.34	0.44
Percent Sulfur	+0.1	+0.10	0.10
Percent Moisture Content	53.0	9.4	89.6

2 1 3327

TABLE 7 SUMMARY OF TCLP ANALYSES RESULTS FOR DRUM SOLIDS PIER DRUM SITE BRADENTON, FLORIDA JANUARY 1995

SAMPLE NUMBER	00911		009	15	00916			
SAMPLE LOCATION	REAC# 1,8,	9.11	REAC	C#5	REAC#3,7,10,13,14			
	CONC	MDL	CONC	MDL	CONC	MDL		
PARAMETER	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L		
BNA's								
Benzyl Alchohol	1000	20	83	20	150	20		
Benzoic Acid	120	20	270	20	4800E	20		
Di-n-butyl phthalate	40	20	34	20	91	20		
phenol	110	20	17J	20	34	20		
2-Methylphenol	860	20	ND	20	ND	20		
METALS								
Cadmium	167	80	ND	4	ND	4		
Calcium	435000	2500	2800	250	1290	250		
Cobalt	470	250	ND	250	ND ND	250		
Copper	ND	50	ND	50	50	50		
Iron	2190	200	ND	200	760	200		
Lead	102	20	7.9	5	6.1	5		
Magnesium	4300	500	300	50	170	50		
Manganese	350	100	ND	100	ND	100		
Potassium	4200	1000	1580	100	570	100		
Sodium	3100000	500000	2300000	500000	1800000	500000		
Zinc	440	50	210	50	130	50		
PEST/PCB's								
BHC-delta	0.6	0.18	ND	0.11	ND	0.25		
Endosulfan-alpha	ND	0.18	ND	0.11	1.0	0.25		
VOC's		<u> </u>				 ,		
Acetone	1300000D	5000	11000J	5000	ND	2500		
Styrene	6100	5000	130000	2000	40000	2500		

MDL = Method Detection Limit

ND = Not Detected

E = Exceeds calibration range, estimated value

J = Detected Below Detection Limit

2 2 3003

TABLE 8 SUMMARY OF GROUNDWATER WELL SAMPLE RESULTS PIER DRUM SITE BRADENTON, FLORIDA JANUARY 1895

SAMPLE NUMBER		00636 MW-S-DUP		00641 MW-N		00835		00634		00637		
SAMPLE LOCATION						MW	-S	MW-W		MW-E		
	MCL	CONC	MDL	CONC	MDL	CONC	MDL	CONC	MDL	CONC	MDL	
PARAMETER	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L	
BNA's												
Di-n-butylphthalate	NEL	1J,B	10.0	50B	11.0	2J,B	10.0	66B	11.0	1J,B	11.0	
METALS (UNFILTERED)												
Aluminum	NEL	1200	40.0	16000	40	1400	40	7400	40	10000	40	
Arsenic	50	3.5	2.2	12.0	2.2	3.0	2.2	10.0	2.2	19.0	2.2	
Barium	2000	18	4.0	44.0	4.0	18.0	4.0	30.0	4.0	39.0	4.0	
Calcium	NEL	4000	100	2300	100	4100	100	27000	100	5800	100	
Chromium	100	ND	2.8	16.0	2.8	ND	2.8	7.6	2.8	13.0	2.8	
Iron	NEL	9400	10	4100	10	9200	10	6600	10	4500	10	
Lead	15	ND	2.2	3.0	2.2	ND	2.2	ND	2.2	3.2	2.2	
Magnesium	NEL	4200	500	1400	500	4300	500	13000	500	5800	500	
Manganese	NEL	4	2.0	5.0	2.0	3.0	2.0	6.0	2.0	8.0	2.0	
Potassium	NEL	ND	2000	ND	2000	ND	2000	2100	2000	ND	2000	
Sodium	*160000	29000	500	3800	500	30000	500	75000	500	40000	500	
Vanadium	NEL	ND	5.0	25.0	5.0	ND	5.0	12.0	5.0	11.0	5.0	
Zinc	NEL	ND	5.0	ND	5.0	ND	5.0	ND	5.0	ND	5.0	
					<u> </u>							
METALS (FILTERED)												
Aluminum	NEL	1000	40	3400	40	1100	40	3500	40	4100	40	
Arsenic	50	3.3	2.2	10.0	2.2	2.8	2.2	10.0	2.2	15.0	2.2	
Barium	2000	17	4	4.0	4.0	17.0	4.0	10.0	4.0	9.0	4.0	
Calcium	NEL	4000	100	1700	100	4100	100	28000	100	6000	100	
Chromium	100	ND	2.8	ND	2.8	ND	2.8	3.8	2.8	3.5	2.8	
iron	NEL	9100	10	1700	10	9200	10	5800	10	3600	10	
Lead	15	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2	
Magnesium	NEL	4100	500	870	500	4200	500	13000	500	5600	500	
Manganese	NEL	3	2.0	ND	2.0	3.0	2.0	3.0	2.0	5.0	2.0	
Potassium	NEL	ND	2000	ND	2000	ND	2000	ND	2000	ND	2000	
Sodium	*160000	29000	500	3000	500	30000	500	76000	500		500	
Vanadium	NEL	ND	5.0	9.0	5.0	ND	5.0	7.0	5.0		5.0	
Zinc	NEL	ND	5.0	ND	5.0	ND	5.0	5.0	5.0	ND	5.0	
PEST/PCB's	. NA	No Analy	/tes Were	Detected	l,							
VOC's												
			2.0	0.4		7.5	2.0	3.7	2.0	8.0		
Acetone	NEI	62	1 211) on	1 / 2	<i></i>				1 211	
Acetone Methyl-tertiary-butylether	NEL NEL	6.2 ND	1.0	3.4 14.0	1.0	ND	1.0	ND	1.0	ND	2.0	

MCL = Maximum Contaminant Level

NOTE: All established MCLs reported indicate Florida State and Federal Regulations.

NEL = No Established MCL

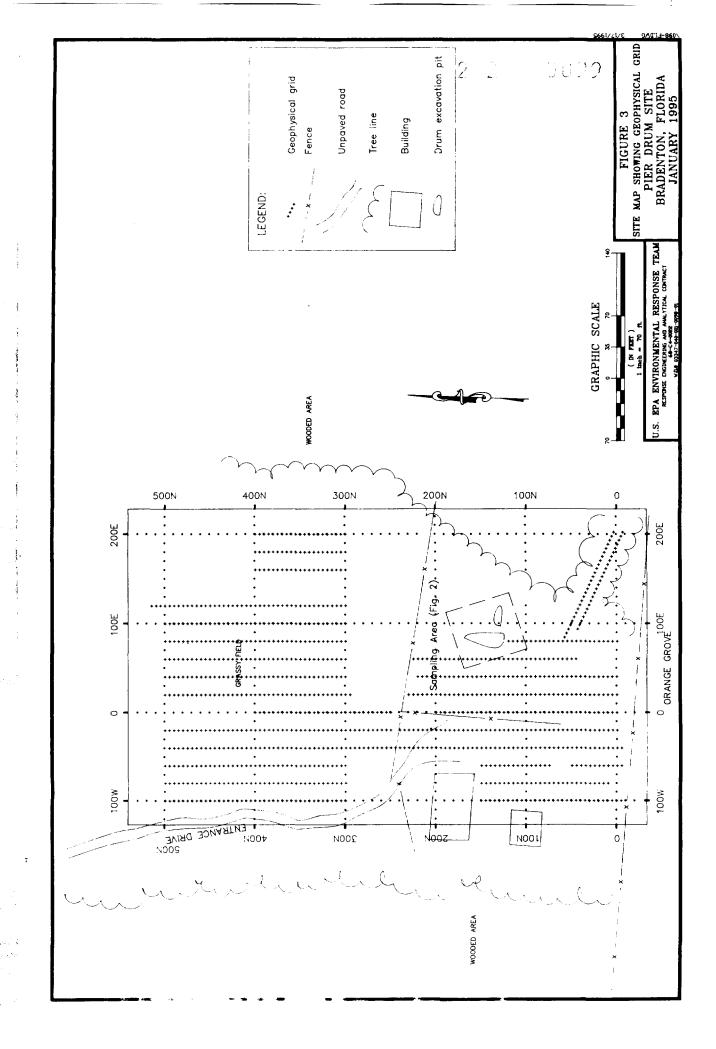
*Florida State Department of Protection MCL

MDL = Method Detection Limit

ND = Not Detected

B = Compound found in blank

J = Detected below detection limit



2 1 9.00

TABLE 9 SUMMARY OF SURFACE WATER SAMPLE RESULTS PIER DRUM SITE BRADENTON, FLORIDA JANUARY 1895

SAMPLE NUMBER		00639 W-PIT		00638 N-PIT	
SAMPLE LOCATION					
	MCL	CONC	MDL	CONC	MDL
PARAMETER	μg/L	μg/L	μg/L	μg/L	μg/L

BNA's					
Di-n-butylphthalate	NEL	47B	10	38B	11

METALS (UNFILTERED))		 		
Aluminum	NEL	2000	40	2300	40
Arsenic	50	10.0	2.2	30	2.2
Barium	2000	9.0	4.0	20	4
Calcium	NEL	12000	100	17000	100
Chromium	100	3.6	2.8	8.6	2.8
Iron	NEL	1600	10	950	10
Lead	15	ND	2.2	ND	2.2
Magnesium	NEL	4600	500	6500	500
Manganese	NEL	26.0	2.0	18	2.0
Potassium	NEL	2300	2000	4900	2000
Sodium	*160000	30000	500	25000	500
Vanadium	NEL	8.0	5.0	7	5
Zinc	NEL	56.0	5.0	27	5

METALS (FILTERED	D)	-			
Aluminum	NEL	1400	40	870	40
Arsenic	50	9.6	2.2	29	2.2
Barium	2000	7.0	4.0	10	4
Calcium	NEL	9800	100	16000	100
Chromium	100	2.9	2.8	4.5	2.8
Iron	NEL	1100	10	560	10
Lead	15	ND	2.2	ND	2.2
Magnesium	NEL	3600	500	5800	500
Manganese	NEL	21.0	2.0	13	2
Potassium	NEL	ND	2000	4400	2000
Sodium	*160000	23000	500	23000	500
Vanadium	NEL	6.0	5.0	ND	5.0
Zinc	NEL	45.0	5.0	25	5.0

PEST/PCB's	NA	No Analytes	Were Detected.

VOC's					
Acetone	NEL	7.3	2.0	ND	2

MCL = Maximum Contaminant Level

NOTE: All MCL reported are for Florida State and Federal Regulations unless otherwise noted.

NEL = No Established MCL

*Florida State Department of Protection MC_

MDL = Method Detection Limit

ND = Not Detected

B = Compound found in blank

J = Detected below detectic limit

TABLE 10 SUMMARY OF RESULTS FOR SEDIMENT SAMPLES PIER DRUM SITE

BRADENTON, FLORIDA JANUARY 1995 2 3101

SAMPLE NUMBER	A00642		A00643		A00644					
SAMPLE LOCATION	SEC)-S	SED	-N	SED	-E				
	CONC	MDL	CONC	MDL	CONC	MDL				
PARAMETER	μg/Kg	μg/Kg	μg/Kg	μg/Kg	μg/Kg	μg/Kg				
BNA's										
Benzyl Alcohol	ND	407	ND	398	ND	402				
Diethylphthalate	87J	407	ND	398	ND	402				
Di-n-butylphthalate	2765B	407	3029B	398	1581B	402				
Butylbenzylphythalate	179J,B	407	156J,B	398	ND	402				
Bis(2-Ethylhexyl)phthalate	338J,B	407	ND	398	39J,B	402				
METALS										
Aluminum	660000	9400	470000	8900	330000	9300				
Barium	4700	3800	3600	3600	ND	3700				
Calcium	290000	47000	94000	45000	61000	47000				
Chromium	1400	800	ND	700	ND	700				
Iron	200000	8400	150000	8000	4300	8400				
Sodium	69000	47000	63000	45000	93000	47000				
Zinc	5400	1900	ND	1800	3400	1900				
PEST/PCB's	No Analyt	es Were D	etected							
VOC's										
Acetone	207	2.5	ND	2.5	ND	2.5				
Methyl-tertiary-butylether	ND	1.3	ND	1.2	ND	1.3				
Toluene	12	1.3	0.9J	1.2	ND	1.3				
Ethylbenzene	670	1.3	ND	1.2	ND	1.3				
o-Xylene	2.4	1.3	ND	1.2	ND	1.3				
Styrene	2.8	1.3	ND	1.2	ND	1.3				
Isopropylbenzene	2.1	1.3	ND	1.2	ND	1.3				
n-Propylbenzene	1.6	1.3	ND	1.2	ND	1.3				
1,2,4-Trimethylbenzene	3.3	1.3	ND	1.2	ND	1.3				

MDL = Method Detection Limit

ND = Not Detected

J = Compound Detected Below Method Detection Limit

2 2 9:02

TABLE 11 SUMMARY OF TCLP ANALYSES RESULTS IN SEDIMENT SAMPLES PIER DRUM SITE BRADENTON, FLORIDA JANUARY 1995

SAMPLE NUMBER	A00525		A00526		A,D00527	
SAMPLE LOCATION	SED-S		SED-		SED-E	
<u> </u>	Conc	MDL	Conc	MDL	Conc	MDL
PARAMETER	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L
BNA's						
Benzyl Alcohol	16	13	17	11	15J	16
Diethylphthalate	5J	13	2 J	11	ND	16
Di-n-butylphthalate	56B	13	45B	11	91B	16
Butylbenzylphythalate	ND	13	4J,B	11	2J,B	16
Bis(2-Ethylhexyl)phthalate	12J,B	13	11B	11	15J,B	16
METALS			<u> </u>		·	
Arsenic	6.3	2.2	ND	2.2	ND	2.2
Barium	930	4.0	28	4.0	11	4.0
Chromium	8.0	5.0	ND	5.0	ND	5.0
Lead	3.2	2.2	ND	2.2	ND	2.2
PEST/PCB's	No Analytes V	Vere Detecte	d			
VOA's	No Analytes V	Vere Detecte	d .			

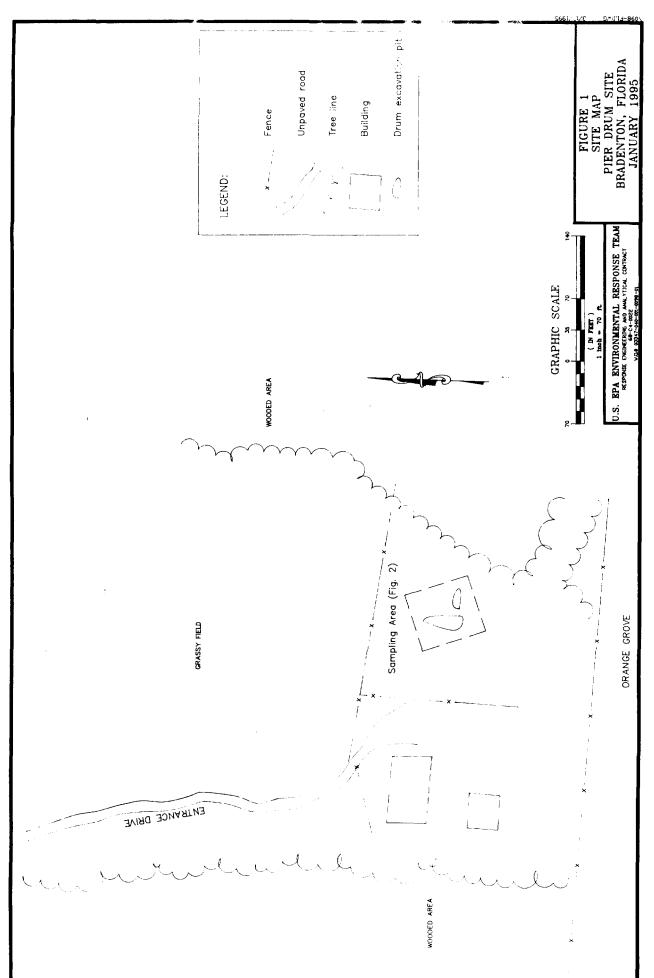
MDL = Method Detection Limit

ND = Not Detected

J = Detected Below Detection Limit

B = Detected in Blank Sample

Figures

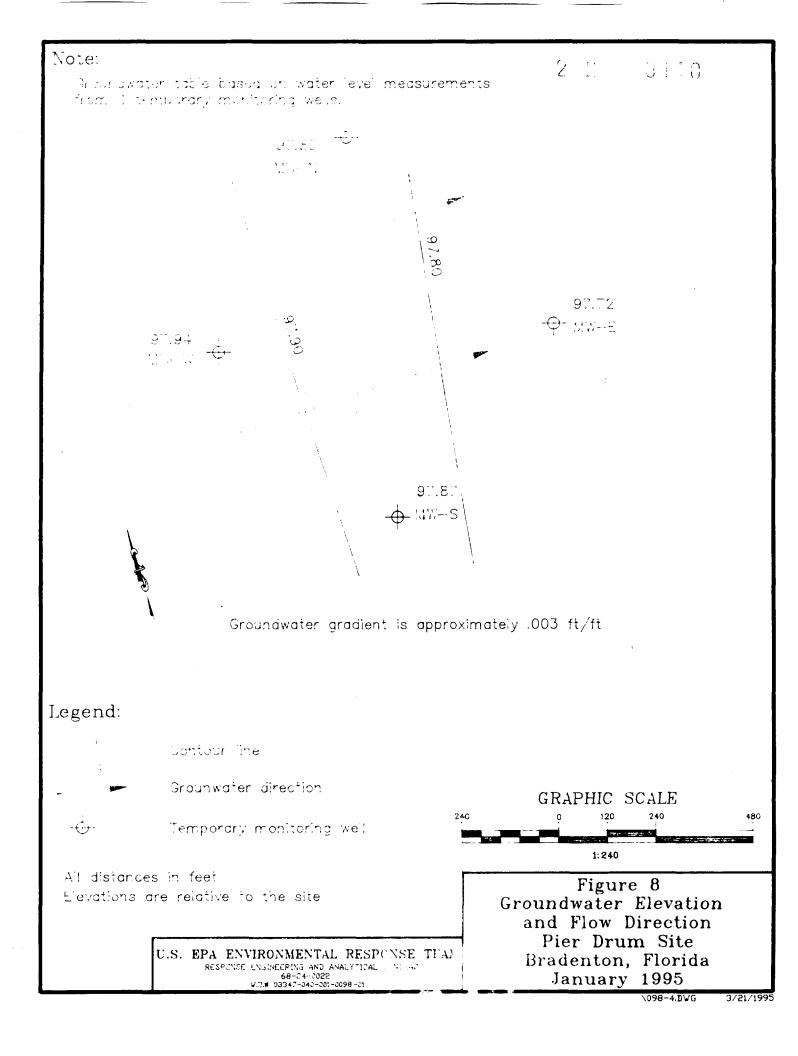


UNSCANNABLE

MEDIA

(PHOTOGRAPHS)

Number of related descent to



Appendix A

APPENDIX A

SAMPLE DOCUMENTATION

Pier Drum Site Bradenton, FL January 1995

Appendix Contents:

Field Notes

Field Data Sheets

Chain of Custudy Communication

.

•

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 2 3 3144

Samplers	no/4. Mengat Drum MW-W	REA	n of Custody No.: C Task Leader: WAM: Assignment No	C. Lewis Powell 0098
SITE DESCRIPTION landfill old field upland palustrin industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	•	SURFACE WATER color odor flow direction	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other sand
groundwater effluent ke groundwater sludge tr potable water leachate b sediment waste a	EVICE emmerer ponar owel other ucket uger kman	SAMPLE INFO color_ odor_ temp_ DO_ cond_	pHORPsalinitysample_depth	
ORGANICS A halogenated & aromatic volatiles B volatiles C trihalomethanes D pesticides/PCB E PCB F base neutral/acid extractables G pesticides drinking water H herbicides drinking water I other INORGANICS A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals other	A total cy B total pl C petrole D. pH E alkalin F hardne G total d H total si I sulfate J. TOC K. grain s L. percen	nenol eum hydrocarbons ty ss ssolved solids uspended solids	SAMPLE PREPA CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVE HNO, NaOH Zn Acetate HCI Na,SO, other
RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS A - Mature F - Mature		H,I- ViDA	´S	

1115

2 %

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

SITE DESCRIPTION Iandfill old feld upland palustrine rock clay color width rock silt industrial wooded lowland riverine grave muck odor depth rubbile clay rubbile clay color muck and remained lacustrine shand loam flow velocity_cm/s grave organic residential gully silt peat direction pools % shell other residential gully silt peat direction pools % shell other residential gully silt peat direction pools % shell other residential gully silt peat direction pools % shell other residential gully silt peat direction pools % shell other residential gully silt peat direction pools % shell other makes floodplain color pH ambient temp color pH ambient temp gully flowers are color pH ambient temp saintly relative humdity sediment waste auger and there odor of ORP barometric pressure residential gully sediment waste auger and the stage and there was the stage of the grave o			~	010	/ · · · · · · ·	Chain of Custody No.:	
Time 795.) Sample Location #GU_S STE DESCRIPTION landfill oid feld upland palustrine rock clay color width rock still motistrail wooded low-land riverine grave muck odor depth ribble clay velocity_cm/s gravel organic residential guily sill peat direction pools \$ shell other older infles \$ sand of the process of the pro	1.10			Kileun)	5 Gettly	REAC Task Leader:	Cenis
Time 795.) Sample Location #GU_S STE DESCRIPTION landfill oid feld upland palustrine rock clay color width rock still motistrail wooded low-land riverine grave muck odor depth ribble clay velocity_cm/s gravel organic residential guily sill peat direction pools \$ shell other older infles \$ sand of the process of the pro	Date//_/	Site Name:	Pier De	erm .	U	EPA WAM: G. PO	welf
STEDESCRIPTION Iandfill old feld upland palustrine rock clay color width rock slit undustral wooded lowland riverine gravel muck odor depth rubble clay rubble clay residential december of the residential december of the residential depth rubble clay rubble clay residential december of the residential dully sit peat direction pools % shell other residential gully sit peat direction pools % shell other residential gully sit peat direction pools % shell other residential gully sit peat direction pools % shell other residential gully sit peat direction pools % shell other residential gully sit peat direction pools % shell other pools % shell other pools % sand residential gully sit peat direction pools % shell other pools % sand residential gully peat	Time			_ S'	•		1998
landfill old field upland palustrine rock clay color depth rock silt rock clay color depth rubble clay clay rubble clay clay rubble clay clay commercia farmland lacustrine sand loam rock clay corner clay color depth rubble clay clay commercia color depth rubble clay corner clay color color clay							
industrial wooded lovaland riverine gravel muck odor depth rubble day commercial familiand lacustrine sand loam flow velocity crivls gravel organic cesidential gully silt peal direction pools % shell other flessedential gully silt peal direction pools % shell other moderate fiftes who sand surface water effluent kemmerer ponar color ph ambient temp broadewater sludge trowel other odor ORP barometric pressure poliate water leachate bucket temp salinity relative humidity seatment waste auger DO sample depth weather conditions weather conditions water leachate bucket temp salinity relative humidity weather conditions water leachate bucket temp salinity relative humidity weather conditions water leachate bucket temp salinity relative humidity weather conditions. ANALYSES TO BE PERFORMED ORGANICS A halogenated & aromatic volatiles A total cynanide glass jar HNO, DRGANICS A halogenated & aromatic volatiles B total phenol plastic part NaOH C trinsionethanes C petroleum hydrocarbons actate core Zn Acetate D pH plastic bucket Na,SO, D pesticides/PCB D pH plastic bucket Na,SO, The recentral placed extractables F hardness C goesticides dirinking water other	SITE DESCRI	PTION		SOIL TYPE	SURFACE WATER	R STREAM	BOTTOM
commercial farmand lacustrine sand loam flow	landfill d	old field upland palus	trine	rock clay			rock silt
residential gully suit peat direction pools % shell other riffles % sand SAMPLE TYPE DEVICE SAMPLE INFORMATION WEATHER PARAMETERS windows floodplain color pH ambient temp many color pH ambient temp salinity relative humidity beatment in each at bucket temp salinity relative humidity sediment waste auger DD sample depth weather conditions. ANALYSES TO BE PERFORMED ORGANICS A halogenated & aromatic volatiles B total phenol plastic par NaCH E volatiles B total phenol plastic par NaCH E PCB E PCB D petroleum hydrocarbons actate core Zn Acatate D pH plastic bag HCl D petrolodes drinking water H total dissolved solids H netrolodes drinking water H total suspended solids A metals, priority pollutant G metals TAL ORGANICS A metals, priority pollutant G metals TAL Ore metals scan (ICP) D metals other Signal FE CACRA A TOLP ORGANICS A metals priority pollutant G metals TAL Ore metals scan (ICP) D metals other Signal FE CACRA A TOLP ORGANICS A metals priority pollutant G metals TAL Ore metals conterned a metals of the signal FE C conswirty DH Oreactivity C corroswirty DH Oreactivity C corroswirty DH Oreactivity C corroswirty DH Oreactivity C corroswirty DH C corroswirty DH Oreactivity C corroswirty DH C CACRA A TOLP C CACRA A TOLP C CACRA A TOLP C CACRA A TOLP C CACRA C C C C C C C C C C C C C C C C C C C	industrial v	wooded lowland rive	rine	gravel muck	odor		rubble clay
SAMPLE TYPE DEVICE SAMPLE INFORMATION WEATHER PARAMETERS	commercial f	farmland lacustrine		sand loam	flow	· ———	gravel organic
SAMPLE TYPE Surface water effluent towel other color pH ambient temp Surface water effluent towel other color pH ambient temp Surface water effluent towel other color pH ambient temp Surface water leachate bucket towel other color pH ambient temp Sediment waste auger DO sample depth weather conditions Social other ekman cond to the stage ANALYSES TO BE PERFORMED ORGANICS A halogenated & aromatic volatiles B volatiles B total phenol plastic par NaOH C trinisiomethanes C petroleum hydrocarbons acetate core Zn Acetate D pesticides/PCB D ph plastic bag HCI E alkalinny plastic bucket Na,SO, other eneutral lacid extractables F hardness other other other C pressionedes drinking water H total suspended solids H netbicides drinking water H total suspended solids C personal representation of the color of	residential g	gully		silt peat	direction		shell other
Surface water effluent kemmerer ponar color PH ambient temp down water sludge trowel other odor ORP barometric pressure potable water leachate bucket temp salinity relative humidity sediment waste auger DO sample depth weather conditions to ther conditions to the stage Soil other ekman Cond It de stage Soil other RANALYSES TO BE PERFORMED ORGANICS OTHER ANALYSES CONTAINER PRESERVATIVES A halogenated & aromatic volatiles B total phenol plastic jar NaOH plastic bag HCI plastic bucket Na,SO, of the salininty plastic bucket Na,SO, of the plastic bucket Na,SO, of the salininty plastic bucket Na,SO, of	hedgerows f	floodplain		color	·	riffles%	sand
Surface water effluent kemmerer ponar color PH ambient temp down water sludge trowel other odor ORP barometric pressure potable water leachate bucket temp salinity relative humidity sediment waste auger DO sample depth weather conditions to ther conditions to the stage Soil other ekman Cond It de stage Soil other RANALYSES TO BE PERFORMED ORGANICS OTHER ANALYSES CONTAINER PRESERVATIVES A halogenated & aromatic volatiles B total phenol plastic jar NaOH plastic bag HCI plastic bucket Na,SO, of the salininty plastic bucket Na,SO, of the plastic bucket Na,SO, of the salininty plastic bucket Na,SO, of	SAMPLE TYPE		DEVICE		SAMPLE	INFORMATION	WEATHER PARAMETERS
Studge S	surface water	effluent	kemmerer	ponar			ambient temp
Dolable water leachate bucket temp salinity relative humidity sesdiment waste auger DO sample depth wather conditions sesdiment waste auger DO sample depth weather conditions sesdiment depth of the sample depth waste stage. ANALYSES TO BE PERFORMED OTHER ANALYSES CONTAINER PRESERVATIVES A halogenated & aromatic volatiles A total cynanide glass jar HNO, plastic jar NaOH D plastic par NaOH D plastic par NaOH D plastic bag HCI D pH plastic bag HCI D p plastic bag HCI E akalinity plastic bucket Na,SO, other oth	groundwater	sludge	trowei	•			
sediment other ekman DO sample depth weather conditions of the destage washer conditions of the destage of the destag	potable water	-					
ANALYSES TO BE PERFORMED ORGANICS A halogenated & aromatic volatiles B volatiles C trihaiomethanes C c petroleum hydrocarbons D peth plastic par D peth plastic bucket D ph D pethodes drinking water H horbicides drinking water H horbicides drinking water C other D TOC C K grain size C MRGANICS A metals priority pollutant C metals scan (ICP) D metals, other O reactivity D comments C C MRGANICS A TOLP C C COTOSWITY D PH D C COTOSWITY D PH D C C C C C C C C C C C C C C C C C C C	sediment	waste					
ORGANICS A halogenated & aromatic volatiles A total cynanide B total phenol D plastic par NaOH C trihalomethanes C petroleum hydrocarbons D pH D plastic bag D petroleum hydrocarbons D plastic bag D pH D plastic par	soil		_				
A halogenated & aromatic volatiles B volatiles B volatiles B total phenol plastic par NaOH C trihalomethanes C petroleum hydrocarbons D pesticides/PCB D pH plastic bag HCI B PCB E alkalinity plastic bucket Na,SO, O ther O particles drinking water O total dissolved solids O total dissolved solids O total dissolved solids O total dissolved solids O total suspended solids O total dissolved solids O total dissolved solids O total dissolved solids O total dissolved solids O total suspended solids O ther O metals priority pollutant O metals priority pollutant O metals scan (ICP) O corrosivity O corros	ANALYSES TO	D BE PERFORMED	 -			SAMPLE PREPAR	ZATION
A halogenated & aromatic volatiles B volatiles B volatiles B total phenol plastic par NaOH C trihalomethanes C petroleum hydrocarbons D pesticides/PCB D pH plastic bag HCI B PCB E alkalinity plastic bucket Na,SO, O ther O particles drinking water O total dissolved solids O total dissolved solids O total dissolved solids O total dissolved solids O total suspended solids O total dissolved solids O total dissolved solids O total dissolved solids O total dissolved solids O total suspended solids O ther O metals priority pollutant O metals priority pollutant O metals scan (ICP) O corrosivity O corros						_	
B volatiles C trihalomethanes C petroleum hydrocarbons acetate core Zn Acetate D pesticides/PCB D pH plastic bag HCI Desticides/PCB E a likalimity plastic bucket Na,SO, other	ORGANICS						
C trihalomethanes C petroleum hydrocarbons D pesticides/PCB D pH plastic bag HCI E alkalinty plastic bucket Na,SO, other Other Other Other I sulfate NORGANICS A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals, other CCOMMENTS A FUNCAL COMMENTS C petroleum hydrocarbons acetate core Zn Acetate HCI E alkalinty plastic bag HCI Na,SO, other Other Other Other I sulfate STORAGE J TOC wet ice dry ice ambient M other Other CCOMMENTS A FUNCAL C corrosvity pH D reactivity other COMMENTS A FUNCAL C COMMENTS A FUNCAL C COMMENTS A FUNCAL C COMMENTS C FUNCAL C C CAP C C C CAP C C C CAP C C C CAP C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C C	-	d & aromatic volatiles		A. total	cynanide	glass jar	•
D pesticides/PCB E PCB E alkalimity D plastic bag D pH D plastic bag D pl D plastic bag D pl D plastic bag D plastic backet D plastic backet D alkalinty D alkalint					•	plastic jar	
E PCB E alkalinity plastic bucket Na,SO, other other Operations destractables F hardness Other O					oleum hydrocarbons	acetate core	
F base neutral/acid extractables G pesticides drinking water H herbicides drinking water H total suspended solids I sulfate J TOC Wet ice NORGANICS A metals, priority pollutant C metals scan (ICP) D metals, other PH O reactivity T other COMMENTS A F Numb E - N	D. pesticides/P	PCB		D. pH		plastic bag	
G pesticides drinking water H herbicides drinking water H total suspended solids I sulfate STORAGE J TOC wet ice NORGANICS K grain size dry ice A metals, priority pollutant B metals. TAL C metals scan (ICP) D metals, other C corrosivity D reactivity E other COMMENTS A FUNCA E GRA E GR	E PCB			E. alkal	linity	plastic bucket	
Herbicides drinking water I sulfate J TOC Wet ice NORGANICS A metals, priority pollutant B metals TAL C metals scan (ICP) D metals other RICRA A TCLP B ignitability C corrosivity D reactivity E other COMMENTS A FLITAL C Metals C	F base neutral	l/acid extractables		F hard	ness	other	
I sulfate STORAGE J TOC wet ice NORGANICS K grain size dry ice A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals other RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS A - Punch F - Pu	G. pesticides, o	drinking water		G total	dissolved solids		
NORGANICS A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals other RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS A - Fund F - Fu	H herbicides of	drinking water		H. total	suspended solids		
NORGANICS A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals, other RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS A - Narah E - Narah	I other			1. sulfat	te	STORAGE	
A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals, other RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS A - Narah F - Narah F - CNA E - CNA						wet ice	
B metals, TAL M other C metals scan (ICP) D metals, other RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS A - March E - GAR E - GAR E - GAR E - GAR	INORGANICS			K grain	n size	dry ice	
Cometals scan (ICP) Dimetals, other RCRA A TCLP Bignitability CorrosivitypH Direactivity E other COMMENTS A - Kinha E -	A metals, prior	rity pollutant				ambient	
D metals, other RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS A - Lide E - Lide T - CAP E - CAP E - CAP E - CAP	B metals, TAL			M othe	er		
RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS $ A = \mu_{stral} F_{stral} $ $ C = \mu_{stral} F_{stral} $							
A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS A - Marak E - Marak (F) C - ENP E - FENP (RE)	U metais, othe	·					
B ignitability C corrosivitypH D reactivity E other COMMENTS A - Funce \[\begin{align*}	RCRA						
C corrosivitypH D reactivity E other COMMENTS A - Name E - Name C - GNA D - GNA E - First felt	A TCLP						
D reactivity E other COMMENTS A - Munde E - Munde C - GNA D - GNA E - First Feb.	B ignitability						
COMMENTS A - Kundi E - Kurel (F) C - BNA D - ENA E - First Feb		pH					
COMMENTS A-FLAND E-FLAND C-BNA D-ENA E-FLAND	D reactivity						
E-firel(F) C-6NA D-6NA E-fire/Feb	E other						
E-firel(F) C-6NA D-6NA E-fire/feb	COMMENTS	1 - 11.300		···=		·	
C-6NA D-ENA E-finitel	COMMENTS		,				
D-ENA E-finifel		= '	•				
E-fear/fel							
E-feer/fel							
CONAL ALL		E- front Re					
Fig. Governor Commission Commissi	FORM #1	F - 631-1968) 2				
9.4 I - 2001		G 4 D - 10	à 15				

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 2 0116

Date 1/11/85 Samplers: 1. Site Name: 15:09 Sample Location	Wir Drum		Chain of Custody No.: REAC Task Leader: EPA WAM: Work Assignment No.:	· Leino · ell O098
SITE DESCRIPTION landfill old field upland palustri industrial wooded lowland riverin commercial farmland lacustrine residential gully hedgerows floodplain	e gravel i sand	PE SURFACE WATER clay color muck odor loam flow peat direction	- · · · · - · · · · · ·	BOTTOM rock silt rubble clay gravel organic shell other sand
surface water effluent groundwater sludge potable water leachate sediment waste	DEVICE kemmerer pona trowel othe bucket auger ekman		ORPsalinitysample depthtide stage	barometric pressure relative humidity weather conditions
ORGANICS A halogenated & aromatic volatiles 3 volatiles C trihalomethanes D pesticides/PCB E PCB F base neutral/acid extractables G pesticides drinking water H herbicides drinking water I other INORGANICS A metals priority pollutant B metals TAL C metals scan (ICP) D metals other		OTHER ANALYSES A. total cynanide B. total phenol C. petroleum hydrocarbons D. pH E. alkalinity F. hardness G. total dissolved solids H. total suspended solids I. sulfate J. TOC K. grain size L. percent moisture M. other	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS: f redxis f redxis	(F)			
C, C - B/H E F - POS G, H, I - V	:/aes			

FIELD DATA SHEET

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 1 3117

Date: 1/13/15 Site Name: Pier Sen Time 854 Sample Location: ML	υ <u>É</u>		Chain of Custody No.: REAC Task Leader: EPA WAM: Work Assignment No.:	
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	SOIL TYPE rock clay gravel muck sand loam silt peat color	SURFACE WATER color odor flow direction	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other
SAMPLE TYPE DEVICE surface water effluent kemmer groundwater sludge trowel potable water leachate bucket sediment waste auger soil other ekman		SAMPLE color odor temp DO cond	ORPsalinitysample_depth	
ANALYSES TO BE PERFORMED ORGANICS A halogenated & aromatic volatiles B volatiles C trihalomethanes D pesticides/PCB E PCB F base neutral/acid extractables G pesticides, drinking water H herbicides, drinking water I other	A. total of B. total of C. petro D. pH E. alkalin F. hardn G. total of total of total of L. sulfate J. TOC	leum hydrocarbons nity ness dissolved solids suspended solids	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
INORGANICS A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals, other RCRA A TCLP B ignitability C corrosivitypH D reactivity E other	•	size int moisture	dry ice ambient	

COMMENTS: A - prefixes

E Metrics: F

CD - GNF

EF - RS+/PCS

FORM#1 GHI - VDFS

REAC, EDISON, NJ (908) 321-4200 **EPA CONTRACT 68-C4-0022**

2 1 7113

Date 1/12/95 Site Name: Fur Sample Location. NF	Drum Pit	Chain of Custody No.:	Verno wall 0-078
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	SOIL TYPE SURFACE WATER rock clay color gravel muck odor sand loam flow silt peat direction color		BOTTOM rock silt rubble clay gravel organic shell other
SAMPLE TYPE surface water effluent kemmer groundwater sludge trowel potable water leachate bucket sediment waste auger soil other ekman	•	ORPsalinitysample_depth	
ORGANICS A halogenated & aromatic volatiles B volatiles C trihalomethanes D pesticides/PCB E PCB F base neutral/acid extractables G pesticides drinking water H herbicides drinking water I other INORGANICS A metals priority pollutant B metals, TAL C metals scan (ICP) D metals other	OTHER ANALYSES A total cynanide B total phenol C. petroleum hydrocarbons D. pH E. alkalinity F hardness G. total dissolved solids H total suspended solids I. sulfate J. TOC K. grain size L. percent moisture M other	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS A - Matals A tratal/F (D - GIVA	· 		
C.D-BIVA E.F- GCG/Fest FORM#1 G.A.I- VOAE			

FIELD DATA SHEET

REAC, EDISON, NJ (908) 321-4200 **EPA CONTRACT 68-C4-0022**

0119

Date 1/11/95 Site Name: Rev 1 Time Sample Location: \(\) (\$\frac{1}{2} \)	Inm STYIT (WPIT)	Chain of Custody No.: REAC Task Leader: EPA WAM: Work Assignment No.:	ox Lewis
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	SOIL TYPE SURFACE WATER rock clay color gravel muck odor sand loam flow silt peat direction color	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other sand
SAMPLE TYPE DEVICE surface water effluent kemmere groundwater sludge trowel potable water leachate bucket sediment waste auger soil otherekman		ORPsalinitysample_depth	barometric pressure relative humidity weather conditions
ANALYSES TO BE PERFORMED ORGANICS A halogenated & aromatic volatiles 3 volatiles C trihalomethanes D pesticides/PCB E PCB F base neutral/acid extractables G pesticides, drinking water H, herbicides, drinking water I other INORGANICS A metals, priority pollutant B metals, TAL C metals, other Indicate the performance of the performan	OTHER ANALYSES A. total cynanide B. total phenol C. petroleum hydrocarbons D. pH E. alkalinity F. hardness G. total dissolved solids H. total suspended solids I. sulfate J. TOC K. grain size L. percent moisture M. other	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
RCRA A TCLP B ignitability C corrosivitypH D reactivity E other			

COMMENTS LE LLO FALL

CD - ENA

FF - Rest / FCE

FORM#1 3 + = Joh 3

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022 2 2 01.0

Date: 1/11/95 Site Name: Sample Location: Fig.	ewis Drin Ell Blonk		Chain of Custody No.: REAC Task Leader: EPA WAM: Work Assignment No.:	Pay Lewis regg Arrell
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	SOIL TYPE rock clay gravel muck sand loam silt peat color	SURFACE WATER color odor flow direction	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other
SAMPLE TYPE surface water effluent kemmer groundwater sludge trowel potable water leachate bucket sediment waste auger soil otherekman		SAMPLE I color odor temp DO cond	ORPsalinitysample_depth	barometric pressure relative humidity weather conditions
ORGANICS A. halogenated & aromatic volatiles B. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water I. other	A. total c B. total p C. petrol D. pH E. alkalii F. hardn G. total c H. total s	ohenol leum hydrocarbons nity ess dissolved solids suspended solids	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
INORGANICS A. metals, priority pollutant B. metals, TAL C. metals scan (ICP) D. metals, other	•	size nt moisture	wet ice dry ice ambient	
A TCLP B ignitability C. corrosivitypH D. reactivity E. other COMMENTS: ### A TCLP ###	.)			

FORM #1

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 2 3121

Date: 1/11/95 Site Name: Time Sample Location:	Pier Druh MW-N		Chain of Custody No.: REAC Task Leader: EPA WAM:Creq Work Assignment No.:	Ray Guis Pausell
SITE DESCRIPTION landfill old field upland palust industrial wooded lowland rivers commercial farmland lacustrine residential gully hedgerows floodplain		y color ck odor am flow	width	BOTTOM rock silt rubble clay gravel organic shell other sand
SAMPLE TYPE surface water effluent groundwater sludge potable water leachate sediment waste soil other	DEVICE kemmerer ponar trowel other_ bucket auger ekman	color	ORPsalinitysample_depth	
ANALYSES TO BE PERFORMED ORGANICS A. halogenated & aromatic volatiles 3. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water I. other INORGANICS A. metals, priority pollutant B. metals, TAL C. metals scan (ICP)	A B C D E F G H I J K L	total cynanide total phenol petroleum hydrocarbons pH alkalinity hardness total dissolved solids total suspended solids sulfate TOC grain size percent moisture	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
D metals, other RCRA A TCLP B. ignitability C corrosivitypH D reactivity E other	- :/s-}.			

FORM #1

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

7.2.22

Date: 1/1/95	Samplers: Site Name: Sample Location:_	SED SED	- Drup		Chain of Custody No.: REAC Task Leader: EPA WAM: Work Assignment No.:	Lay lewis Gregg Rosell
SITE DESCRIPTION Industrial wood commercial farmly residential gully hedgerows flood	eld upland palust ded lowland riveri land lacustrine	rine roo ne gra sai silt	avel muck nd loam	SURFACE WATER color odor flow direction	STREAM width depth velocity cm/s pools% riffles%	shell other
SAMPLE TYPE surface water groundwater potable water sediment soil	effluent sludge leachate waste other	DEVICE kemmerer trowel bucket auger ekman	ponar other	SAMPLE color odor temp DO cond	ORPsalinitysample_depth_	
ANALYSES TO BE ORGANICS A. halogenated & a 3. volatiles C trihalomethanes D pesticides/PCB E. PCB F. base neutral/acid G pesticides, drink H. herbicides, drink I. other INORGANICS A metals, priority p B metals, TAL C. metals scan (ICF D metals, other RCRA	d extractables ing water ing water		A. total B. total C. petro D. pH E. alkal F. hard G. total H total I sulfat J. TOC K. grain L. perco	ness dissolved solids suspended solids e	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
A TCLP B ignitability C corrosivity D reactivity E other		7CL	- ',' O	ell TCL	f except	VOAS

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 1 1123

Date:	SOIL TYPE ine rock clay	RE/	sin of Custody No.:AC Task Leader: A WAM:COOK K Assignment No.: STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other sand
SAMPLE TYPE surface water effluent groundwater sludge potable water leachate sediment waste soil other	DEVICE kemmerer ponar trowel other bucket auger ekman	SAMPLE INFO color odor temp DO cond	DRMATION pH ORP salinity sample depth tide stage	WEATHER PARAMETERS ambient temp barometric pressure relative humidity_ weather conditions
ORGANICS A. halogenated & aromatic volatiles 3. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water I. other INORGANICS A. metals, priority pollutant B. metals, TAL C. metals, other	A. total of B. total of C. petro D. pH E. alkalin F. hardn G. total of tota	leum hydrocarbons nity less dissolved solids suspended solids	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
RCRA A. TCLP B. ignitability C. corrosivitypH D. reactivity E. other COMMENTS: A	2 TCL & TCL	TLP xee	et Voas	

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 1 3124

Date: //II/95 Site Name: Time: Sample Location:	Pier Drun	RI	PA WAM:	lay Lewis 70 Gregg Powell
SITE DESCRIPTION landfill old field upland palustrin industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	•	SURFACE WATER color odor flow direction	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other
surface water effluent degroundwater sludge to potable water leachate sediment waste as soil other	DEVICE lemmerer ponar rowel other bucket luger lekman	SAMPLE INF color odor temp DO cond	ORPsalinitysample_depthtide_stage	barometric pressure relative humidity weather conditions
ANALYSES TO BE PERFORMED ORGANICS A. halogenated & aromatic volatiles 3 volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water II. other	A. total B. total C. petro D. pH E. alkal F. hardi G. total H. total I. sulfati J. TOC K. grain L. perce	inity ness dissolved solids suspended solids	CONTAINER glass jar plastic jar acetate core plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO ₃ NaOH Zn Acetate HCI Na,SO ₄ other
RCRA A. TCLP B. ignitability C. corrosivitypH D. reactivity E other COMMENTS \$\int \int \text{ \text{ \int \text{ \text{ \text{ \text{ \te	•	if excep	7 VOG5	

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 2 9125

SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully	SOIL TYPE SURFACE WATER e rock clay color gravel muck odor sand loam flow silt peat direction	Chain of Custody No.: REAC Task Leader: PA WAM: STREAM Width rock rubble velocity cm/s gravel organic pools % shell other
surface water effluent ke groundwater sludge tro potable water leachate bu sediment waste au	color EVICE SAMPLE IN color emmerer ponar color owel other odor ucket temp uger DO kman cond	ORP barometric pressure salinity relative humidity sample depth weather conditions
ANALYSES TO BE PERFORMED ORGANICS A. halogenated & aromatic volatiles 3. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water I. other INORGANICS A. metals, priority pollutant B. metals, TAL C. metals scan (ICP) D. metals, other RCRA	OTHER ANALYSES A. total cynanide B. total phenol C. petroleum hydrocarbons D. pH E. alkalinity F. hardness G. total dissolved solids H. total suspended solids I. sulfate J. TOC K. grain size L. percent moisture M. other	SAMPLE PREPARATION CONTAINER PRESERVATIVES glass jar HNO, plastic jar NaOH acetate core Zn Acetate plastic bag HCI plastic bucket Na,SO, other other STORAGE wet ice dry ice ambient
A TCLP B ignitability C corrosivitypH D reactivity E other		

0094	Drew Sam	alin		
101295				
	REAC#1			
			Poto 11	
-	OI FLAX	DEP#		
-				
-	The Dm is	on its so	le to ver	
-	Downged			
-	Microtip at Bu SOCED - Bro	ing somo	ist I My be	Lecking
<u> </u>	microtip at Bu	ng 126 ppn	115-11	
•	SOUTD - Bro	5W W 3/	4 / 4 / /	
•				
i			·	
1				** - **
:				
•	· ·			
-	· · · · · · · · · · · · · · · · · · ·			#*
•				
•				
·				
			·	
٠.			•	
				_
				
				· · · · · ·

2 1 1127

Samplers: Getty & Morganti Date: 10 Jan 95 Work Assignment Number: 3347-04-01-0098 Site Name: Pier Drum Container Number/Sample Number: REAC#1 REAC Task Leader: Ray Lewis SITE INFORMATION: 1. Terrain, drainage description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's °F and sunny. MET station on site: X No, SAMPLE INFORMATION: 1. Container type: X Drum, Tank, Other: Shape: 17E in overpak 17H 9 2. Container dimensions: Approximate size: 55 gal in 85 gal 3. Label present: X No. Yes: 4. Spill or Leak present: X No, Yes, Dimensions: sludge Locks liquid, X solid (powder or X crystals), 5. Description: 6. Color: Brown , Odor: Yes , Vapors: -Yes-FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING IGNITABILITY (V Flash X No Flash pH OF AQUEOUS SOLUTION_ pH nod the HOU used in Solubility Fort 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATER REACTIVE! Results: _____ total, ____ partial, ____ no solubility. WATER REACTIVITY: 1. Add small amount of sample to water: _____ bubbles, ____ color change to _____, odor or vapor formation, _____ heat, __X_ No Change. SPECIFIC GRAVITY TEST (compared to water): Sinks, Floats. I. Add small amount of sample to water: 2. If liquid sample sinks, screen for chlorinated compounds. If liquid sample floats and appears to be oily, screen for PCB's (Chlor n Oil kit). CHLOR N OIL TEST KIT INFORMATION: 1. Test kit used for this sample: ____ Yes, ____ No PCB not present; PCB present, less than 50 ppm; 2. Results: PCB present, greater than 50 ppm; _____ 100 % PCB present. SPILL-FYTER CHEMICAL CLASSIFIER STRIPS: MA Results:

When the Solid was phased in the Fligh tester a harder for I min it did 1=1254.

+ for bet , Dra, . Tube

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 1 5123

Samplers:		TER STREAM width depth velocity cm/s	/.
surface water effluent kei groundwater sludge tro potable water leachate bur sediment waste au	mmerer ponar color wel other odor cket temp ger DO_	salinitysample_depth	WEATHER PARAMETERS _ambient tempbarometric pressurerelative humidityweather conditions
ORGANICS A. halogenated & aromatic volatiles d. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water i other INORGANICS A. metals, priority pollutant B. metals, TAL C. metals, other	OTHER ANALYSES A. total cynanide B. total phenol C. petroleum hydrocarbor D. pH E. alkalinity F. hardness G. total dissolved solids H. total suspended solids I. sulfate J. TOC K. grain size L. percent moisture M. other	CONTAINER glass jar plastic jar acetate core plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
RCRA A. TCLP B. ignitability C. corrosivitypH D. reactivity E. other COMMENTS Drum Sample FLA DEP#	(REAC#1		

Samplers: Getty & Morganti 10 Jan 95 Date: Work Assignment Number: 3347-04-01-0098 Site Name: Pier Drum Container Number/Sample Number: REAC#2 REAC Task Leader: Ray Lewis SITE INFORMATION: 1. Terrain, drainage description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's F and sunny. MET station on site: X No. Yes SAMPLE INFORMATION: 1. Container type: X Drum, Tank, Other: 2. Container dimensions: Shape: 17E in overpak 17H M Approximate size: 55 gal in 85 gal 3. Label present: X No, Yes: 4. Spill or Leak present: X No, Yes, Dimensions:
5. Description: X liquid, X solid (powder or X crys powder or ** crystals), sludge 6. Color: , Odor: Yes FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING **IGNITABILITY** No Flash X Flash pH OF AQUEOUS SOLUTION 1 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATERREACTIVE!] Results: _____ total, ____ partial, ____ no solubility. WATER REACTIVITY: 1. Add small amount of sample to water: _____ bubbles, ____ color change to ____ odor or vapor formation, ____ heat, ___ No Change. SPECIFIC GRAVITY TEST (compared to water): 1. Add small amount of sample to water: X Sinks, X Floats. 2. If liquid sample sinks, screen for chlorinated compounds. If liquid sample floats and appears to be oily, CHLOR N OIL TEST KIT INFORMATION: - Test dan 2 times.

1. Test kit used for this sample: X Yes, No

2. Results: PCB not present; PCB present, less than 50 ppm;

X PCB present, greater than 50 ppm; 100 % PCB present.

Filed the Quarti Chlor Test (Hemming PCB's are present but Could not spill the Quarti Chlor Test (Hemming PCB's are present but Could not spill the Quartify)

SPILL-FYTER CHEMICAL CLASSIFIER STRIPS: Quartify

Octor > screen for PCB's (Chlor n Oil kit).

Results: Positive for Petrolan Product, Organic Solvant & Chlorine Rick COMMENTS

1) liquid Separts atth Smi is dissolved w/ the. Part Floods a Part Sinks.

3) Chlorin wine test produce is a firm

4) Product sprens to be Sin by a

y me apposed to time.

REAC, EDISON, NJ (908) 321-4200 **EPA CONTRACT 68-C4-0022**

Date: 18 0		Pier On Sera:	4	REAC Task Leader: Rsy Lewis EPA WAM: Grey lowell Work Assignment No.: 33 77 400000000000000000000000000000000		
SITE DESCI landfill industrial commercial residential hedgerows	old field upland palus wooded lowland rive	strine (SOIL TYPE rock clay gravel muck sand loam silt peat color	SURFACE WATER color odor flow direction	width depth velocity cm/s	BOTTOM rock silt rubble clay gravel organic shell other sand
SAMPLE TY surface wate groundwater potable wate sediment soil	er effluent sludge	DEVICE kemmerer trowel bucket auger ekman	ponar other	SAMPLE color odor temp DO cond	ORPsalinitysample_depth	
ANALYSES	TO BE PERFORMED		 		SAMPLE PREPAR	RATION
ORGANICS halogenated & aromatic volatiles by volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water I. other INORGANICS A. metals, priority pollutant B. metals, TAL C. metals, other D. metals, other		OTHER ANALYSES A. total cynanide B. total phenol C. petroleum hydrocarbons D. pH E. alkalinity F. hardness G. total dissolved solids H. total suspended solids I. sulfate J. TOC K. grain size L. percent moisture M other		container glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other	
RCRA A. TCLP B. ignitability C. corrosivity D reactivity E other						

PLA DEP#1 17E

/	
-	RCAC*3
- 	03 FLADER # Bh 15
- -	178
- •	
TRO70EHL LEW	Day is Dented and in Cut open
- معلم ملده	on the Side by bury
07031	on 10 siece by burg
190	7/0 [11 colid 2/och: 10/
C -	7/8 Full solid plastic like Very hard to sample Ambar in color
	Very hard to senfle
	Hmber IN Color
	· · · · · · · · · · · · · · · · · · ·
-	
cialada _{MM}	
\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \	
r.d_	
÷_	
·	
· _	
	•
~	
_	
•	
· , , –	
-	
-	
_	

2 1 0133

10 Jan 95 Samplers: Getty & Morganti Date: Site Name: Pier Drum Work Assignment Number: 3347-04-01-0098 REAC Task Leader: Ray Lewis Container Number/Sample Number: REAC#3 SITE INFORMATION: 1. Terrain, drainage description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's °F and sunny. MET station on site: X No, Yes SAMPLE INFORMATION: 1. Container type: X Drum, Tank, Other: 2. Container dimensions: Shape: 17E in overpak 17H Approximate size: 55 gal in 85 gal 3. Label present: X No, Yes: 4. Spill or Leak present: X No, Yes, Dimensions: 5. Description: liquid, X solid (powder or X crystals), sludge Pink -> Clow, Odor: Yes 6. Color: , Vapors: Yes FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING **IGNITABILITY** No Flash Flash pH OF AQUEOUS SOLUTION 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATERREACTIVE!] Results: _____ total, ____ partial, ____ no solubility. WATER REACTIVITY: 1. Add small amount of sample to water: _____ bubbles, ____ color change to _____ odor or vapor formation, ____ heat, ___ No Change. SPECIFIC GRAVITY TEST (compared to water): 1. Add small amount of sample to water: X Sinks, Floats. 2. If liquid sample sinks, screen for chlorinated compounds. If liquid sample floats and appears to be oily, screen for PCB's (Chlor n Oil kit). CHLOR NOIL TEST KIT INFORMATION:

1. Test kit used for this sample: _____ Yes ____ No

2. Results: ____ PCB not present; ____ PCB present, less than 50 ppm; ____ 100 % PCB present.

+ A. Signer Dragen Take

Results:

COMMENTS

SPILL-FYTER CHEMICAL CLASSIFIER STRIPS:

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022 2 : 0:34

SITE DESCI landfill industrial commercial residential	old field upland p	tion: SZ72.50 palustrine riverine	SOIL TYPE rock clay gravel much sand loan silt peat		EPA WAM:	BOTTOM rock silt rubble clay gravel organic shell other sand
SAMPLE TY		DEVICE		SAMPLE	INFORMATION	WEATHER PARAMETERS
surface wate	er effluent	kemmere	r ponar	color	pH	_ ambient temp
groundwater	sludge	trowel	other	odor	ORP	barometric pressure
potable wate	er leachate	bucket		temp		relative humidity
sediment	waste	auger		DO		weather conditions
soil	other	ekman		cond	tide stage	
ANALYSES	TO BE PERFORMED)			SAMPLE PREPAR	RATION
ORGANICS			от	HER ANALYSES	CONTAINER	PRESERVATIVES
halogenat	ed & aromatic volatile	es	Α.	total cynanide	glass jar	HNO,
B volatiles			В.	total phenol	plastic jar	NaOH
C. trihalomet	···		C.	petroleum hydrocarbons	acetate core	Zn Acetate
D. pesticides	s/PCB		D. pH		plastic bag	HCI
E PCB			E. alkalinity		plastic bucket	Na,SO,
	ral/acid extractables		F hardness		other	other
	s, drinking water			total dissolved solids		
	s drinking water			total suspended solids ulfate	STORAGE	
· other		-		TOC	wet ice	
INORGANIC	S	<u></u>	_	grain size	dry ice	
A. metals, pr	ority pollutant			percent moisture	ambient	
B metals, TA			M.	other		
C. metals sc	an (ICP)					
D. metals, ot	her					
RCRA						
A TCLP						
B. ignitability						
	/pH					
D reactivity						
E. other		_				

COMMENTS

Drum Sample Refe#3
FLA DEP#3
17E

REAC#4 Q4 FLADEP# Salveye 17H	Psto 16
DM 15 not Second Drum in over pack 2/8 Felt Sold place Very hard to s Ambore In C	which ike CP
LIQUID With s 2ppm /4 ful CLEAR Liquid	

2 2 1136

10 Jan 95 Date: Samplers: Getty & Morganti Work Assignment Number: 3347-04-01-0098 Site Name: Pier Drum REAC Task Leader: Ray Lewis Container Number/Sample Number: REAC#4 SITE INFORMATION: 1. Terrain, drainage description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's °F and sunny. MET station on site: X No. SAMPLE INFORMATION: 1. Container type: X Drum, Tank, Other: 2. Container dimensions: Shape: 17E in overpak 17H Approximate size: 55 gal in 85 gal 3. Label present: Yes: X No, 4. Spill or Leak present: X No, Yes, Dimensions: X solid (powder or 5. Description: Y liquid, X-orystals), 6. Color: , Odor: Yes Nb , Vapors: Yes FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING **IGNITABILITY** No Flash Flash pH OF AQUEOUS SOLUTION. 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATERREACTIVE!] Results: _____ total, _____ partial, _____ no solubility. WATER REACTIVITY: 1. Add small amount of sample to water: bubbles, ___ color change to _____ Add small amount of sample to water: bubbles, odor or vapor formation, heat, No Change. SPECIFIC GRAVITY TEST (compared to water): Sinks, 1. Add small amount of sample to water: Floats. 2. If liquid sample sinks, screen for chloringsed compounds. If liquid sample floats and appears to be oily, screen for PCB's (Chlor n Oil kit).

COMMENTS

CHLOR NOIL TEST KIT INFORMATION:

PCB present, greater than 60 ppm; _____ 100 % PCB present.

SPILL-FYTER CHEMICAL CLASSIFIER STRIPS:

PCB present, less than 50 ppm;

Yes,

Seems to be a but sorpy.

1. Test kit used for this sample: ____

2. Results:

Results:

____ PCB not present;

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022 2 2 0 137

Samplers: <u>Getty</u> Date: <u>10 Jan 15</u> Site Name: <u>Rêr 1</u> Time: <u>1300</u> Sample Location: <u>Ser</u>	Chain of Custody No.: 3627 REAC Task Leader: Lay Leurs EPA WAM: Grey Bowell Work Assignment No.: 349400404507			
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	SOIL TYPE SURFACE WATE rock clay color gravel muck odor sand loam flow silt peat direction color	width depth velocity cm/s	BOTTOM rock silt rubble clay gravel organic shell other sand	
SAMPLE TYPE DEVIC surface water effluent kemme groundwater sludge trowel potable water leachate bucket sediment waste auger soil other ekman	other color temp	ORPsalinitysample_depth		
ANALYSES TO BE PERFORMED		SAMPLE PREPAR	RATION	
ORGANICS A. halogenated & aromatic volatiles B. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water I. other INORGANICS A. metals, priority pollutant B. metals, TAL C. metals scan (ICP) D. metals, other	OTHER ANALYSES A. total cynanide B. total phenol C. petroleum hydrocarbons D. pH E. alkalinity F. hardness G. total dissolved solids H. total suspended solids I. sulfate J. TOC K. grain size L. percent moisture M. other	container glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other	
RCRA A TCLP B. ignitability C corrosivitypH D reactivity E. other				

COMMENTS.

Drum Sousphe REACH 4
FLA DEP #4
Solvere Dm 1714

30	2 .2	9+48	
REAC #5	<u></u>		and the second s
05 FLAX Solvey 1	0EP# 1	36 17	
પ્રાપ્ય	<i></i>		
400 ppm mi		FULL	
CLEAR	GELATANA Pack	cious sl	udge
IN OVER	pack		
			· · · · · · · · · · · · · · · · · · ·
		/	
			u
	<u> </u>		
			•
		· · · · · · · · · · · · · · · · · · ·	

Samplers: Getty & Morganti Site Name: Pier Drum Container Number/Sample Number: REAC#5	Date: 10 Jan 95 Work Assignment Number: 3347-04-01-0098 REAC Task Leader: Ray Lewis
SITE INFOR	MATION:
 Terrain, drainage description: Flat, roul woods and Weather conditions (from observation): Mid to high MET station on site: X No, Yes 	· · · · · · · · · · · · · · · · · · ·
SAMPLE INFO	ORMATION:
1. Container type: X Drum, Tank, Other: 2. Container dimensions: Shape: 17E in overpal Approximate size: 55 3. Label present: X No, Yes: 4. Spill or Leak present: X No, Yes, Description: Ilquid, X Solid (powder or 6. Color: Yes)	gal in 85 gal a in 85 gal imensions: X-crystals), sludge , Vapors: Xee M
FIELD TEST DATA SHEET FO	OR DRUM/TANK SAMPLING
IGNITAB Flash	
pH OF AQUEOU 1. Using 0-14 pH paper, check pH of water/sample solu	IS SOLUTION PLA.
WATER SOLUE 1. Add approximately one part sample to five parts wat HEAT IF WATERREACTIVE!] Results: tot	BILITY TEST: ter. You may need to stir and heat gently. [DO NOT
WATER REAL 1. Add small amount of sample to water: bubble odor or vapor formation, heat, N	es, color change to ,
SPECIFIC GRAVITY TE. 1. Add small amount of sample to water: Sinks, 2. If liquid sample sinks, screen for chlorinated compound screen for PCB's (Chlorin Oil kit).	, Floats.

SPILL-FYTER CHEMICAL CLASSIFIER STRIPS:

CHLOR N OIL TEST KIT INFORMATION:

1. Test kit used for this sample: Yes, No

2. Results: PCB not present; PCB present, less than 50 ppm; PCB present, greater than 50 ppm; 100 % PCB present.

Results:

COMMENTS

Seems to be an alhesine + for & Touler Drin-Tube

00605

FIELD DATA SHEET

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022 2 2 0140

Samplers: Getty, Morganti Date: Din 95 Site Name: Pier Druns Time 1300 Sample Location: Samota PC4				Chain of Custody No.: 3627 REAC Task Leader: Ray Lewis EPA WAM: Gray Awell Work Assignment No.: 3547 40010008		
•	palustrine d riverine rine	SOIL TYPE rock clay gravel muck sand loam silt peat color	SURFACE WATER color odor flow direction	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other sand	
SAMPLE TYPE surface water effluent groundwater sludge potable water leachate sediment waste soil other	DEVICE kemmerer trowel bucket auger ekman	ponar other	SAMPLE color odor temp DO cond	ORPsalinitysample_depth	WEATHER PARAMETERS ambient temp barometric pressure relative humidity weather conditions	
ANALYSES TO BE PERFORME ORGANICS A halogenated & aromatic volati B volatiles C trihalomethanes D pesticides/PCB E PCB F base neutral/acid extractables G pesticides, drinking water H herbicides, drinking water I other INORGANICS A metals, priority pollutant	les	A. total B. total C. petro D. pH E. alka F. hard G. total H. total I. sulfal J. TOC K. grain	ness dissolved solids suspended solids ee	SAMPLE PREPAR CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other	
A. metals, priority pollutant B. metals, TAL C. metals scan (ICP) D. metals, other RCRA A. TCLP B. ignitability C. corrosivitypH D. reactivity E. other		•	ent moisture er	ambient		

COMMENTS

Drus Sunple Rete #5

FLA DEP # 5 Selvare Da 1714/

. =	REAC# C		2	a" ►	1144
	DG FLADEP # 17E				
	178				
····					
	Don is in fair Shape	_			
		_			
	Multicolored Solid				
-	Multicolored solid Contents appear to have polymerized NO-SAMPHE				
	NO-SAMPLE				
			,		
-		_ ·			
		- ·			
	· · · · · · · · · · · · · · · · · · ·	1			
		- !			
		1			
		}	ŀ		
		-			
		- 1	-		
-					
	· · · · · · · · · · · · · · · · · · ·	} .			
		-			
		•			
£					
	· · · · · · · · · · · · · · · · · · ·		ľ		

REAC # 7 Rusty Blue + white drum 7/8 full Solid Multicolored NESIN Very hard to sa-ple.

2 2 3143

Samplers: Getty & Morganti Date: 10 Jan 95 Work Assignment Number: 3347-04-01-0098 Site Name: Pier Drum REAC Task Leader: Ray Lewis Container Number/Sample Number: REAC#7 SITE INFORMATION: 1. Terrain, drainage description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's °F and sunny. MET station on site: X No, Yes SAMPLE INFORMATION: Tank, Other: 1. Container type: X Drum, 2. Container dimensions: Shape: 17E in overpak 17H Approximate size: 55 gal in 85 gal 3. Label present: X No, Yes: 4. Spill or Leak present: X No, Yes, Dimensions: X solid (powder or X crystals), 5. Description: liquid, sludge 6. Color: , Odor: Yes , Vapors: Yes-FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING **IGNITABILITY** No Flash Flash pH OF AQUEOUS SOLUTION 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATER REACTIVE!] Results: _____ total, ____ partial, ____ no solubility. WATER REACTIVITY: 1. Add small amount of sample to water: _____ bubbles, ____ color change to _____ odor or vapor formation, ____ heat, ____ No Change. SPECIFIC GRAVITY TEST (compared to water): 1. Add small amount of sample to water: ______ Sinks, _____ Floats. 2. If liquid sample sinks, screen for chlorinated compounds. If liquid sample floats and appears to be oily, screen for PCB's (Chlor n Oil kit). CHLOR N OIL TEST KIT INFORMATION: 1. Test kit used for this sample: Yes, No
2. Results: PCB not present; PCB present, less than 50 ppm; PCB present. SPILL-FYTER CHEMICAL CLASSIFIER STRIPS:

Product Burns

Results:

COMMENTS

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 2 31/4

Samplers: Samplers: Samplers: Samplers: Sample Location:	Servich PU	EPA WAM: Green	Auell Ruell 354 HO40100085
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	e rock clay color gravel muck odor sand loam flow	RFACE WATER STREAM or width or depth velocity cm/s ction pools % riffles %	•
surface water effluent kee groundwater sludge tro potable water feachate bu sediment waste au	EVICE emmerer ponar owel other ucket uger kman		barometric pressure
ANALYSES TO BE PERFORMED ORGANICS A. halogenated & aromatic volatiles B. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water	OTHER ANALY A. total cynanic B. total phenol C. petroleum h D. pH E. alkalinity F. hardness G. total dissolv H. total suspen I. sulfate J. TOC	de glass jar plastic jar sydrocarbons acetate core plastic bag plastic bucket other	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
INORGANICS A. metals, priority pollutant B metals, TAL C. metals scan (ICP) D metals, other RCRA A TCLP	K. grain size L. percent mois M. other	dry ice sture ambient	
B ignitability C corrosivitypH D reactivity E other			

COMMENTS

Drun Sample Rete #7
PLA DEP#7
17E

REK#8
8 FLADER# Photo 18
5 x voge 174
28 02
1/2 Full 3.8 ppm
Brown + white resin sludge With sand
·
·

Date:

10 Jan 95

Samplers: Getty & Morganti Site Name: Pier Drum Work Assignment Number: 3347-04-01-0098 Container Number/Sample Number: REAC#8 REAC Task Leader: Ray Lewis SITE INFORMATION: 1. Terrain, drainage description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's °F and sunny. MET station on site: X No, Yes SAMPLE INFORMATION: 1. Container type: X Drum, Tank, Other: 2. Container dimensions: Shape: 17E in overpak 17H Approximate size: 55 gal in 85 gal 3. Label present: X No, Yes: Yes, Dimensions: 4. Spill or Leak present: X No, 5. Description: liquid, X solid (powder or sludge X crystals), 6. Color: , Odor: Yes , Vapors: Yes FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING **IGNITABILITY** No Flash Flash ph of aqueous solution 7 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATER REACTIVE!] Results: _____ total, ____ no solubility. WATER REACTIVITY: 1. Add small amount of sample to water: ____ bubbles, __ color change to _____ __ odor or vapor formation, ____ heat, ____ No Change. SPECIFIC GRAVITY TEST (compared to water): 1. Add small amount of sample to water: X Sinks, Floats. 2. If liquid sample sinks, screen for chlorinated compounds. If liquid sample floats and appears to be oily, screen for PCB's (Chlor n Oil kit). CHLOR NOIL ZEST KIT INFORMATION: 1. Test kit used for this sample: Yes, No

2. Results: PCB not present; PCB present, less than 50 ppm; PCB present. PCB present.

COMMENTS

SPILL-FYTER CHEMICAL CLASSIFIER STRIPS:

Results:

00608

2 2 11 17

REAC, EDISON, NJ (908) 321-4200 **EPA CONTRACT 68-C4-0022**

Samplers:	Mazzaki	Chain of Custody No.:	Levis
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	SOIL TYPE rock clay color gravel muck odor sand loam flow silt peat direction color	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other
SAMPLE TYPE DEVICE surface water effluent kemmere groundwater sludge trowel potable water leachate bucket sediment waste auger soil otherekman		ORPsalinitysample_depth	WEATHER PARAMETERS ambient temp_ barometric pressure_ relative humidity_ weather conditions
ANALYSES TO BE PERFORMED ORGANICS halogenated & aromatic volatiles B. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water	OTHER ANALYSES A. total cynanide B. total phenol C. petroleum hydrocarbons D. pH E. alkalinity F. hardness G. total dissolved solids H. total suspended solids I. sulfate	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
INORGANICS A. metals, priority pollutant B. metals, TAL C. metals scan (ICP) D. metals, other	J. TOC K. grain size L. percent moisture M. other	wet ice dry ice ambient	
A TCLP B. ignitability C. corrosivitypH D. reactivity E. other			

Drum Sample REAC#8

FLA DEP #8

Selvase Dm 1714/

FORM #1

REAC #9

BY PLABER # Parts 19

Solvere 1714

Drum appside down in overpack ColLiquid in batt 6P

Drum covered with send bung

Obscured.

Liquid in Bottom of overpack

1000 ppm on micro pip.

DRUM 1/2 Full

1/2 - liquid - Cleer

1/2 - Slage/resin - gray Brown

Date:

10 Jan 95

Samplers: Getty & Morganti Work Assignment Number: 3347-04-01-0098 Site Name: Pier Drum Container Number/Sample Number: REAC#9 REAC Task Leader: Ray Lewis SITE INFORMATION: 1. Terrain, drainate description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's F and sunny. MET station on site: X No, Yes SAMPLE INFORMATION: 1. Container type: X Drum, Tank, Other: 2. Container dimensions: Shape: 17E in overpak 17H Approximate size: 55 gal in 85 gal 3. Label present: X No. Yes: Yes, Dimensions: 4. Spill or Leak present: X No, solid (
Odor: Yes 5. Description: X liquid, X solid (powder or X crystals), sludge 6. Color: led clay , Vapors: Yes FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING **IGNITABILITY** No Flash Flash pH OF AQUEOUS SOLUTION 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATERREACTIVE!] Results: _____ total, ____ partial, ____ no solubility. WATER REACTIVITY: 1. Add small amount of sample to water: ____ bubbles, X color change to Cloudy Very odor or vapor formation, heat, No Change. SPECIFIC GRAVITY TEST (compared to water): Sinks, 1. Add small amount of sample to water: Floats. 2. If liquid sample sinks, screen for chlorinated compounds. If liquid sample floats and appears to be oily, screen for PCB's (Chlor n Oil kit). CHLOR N OIL TEST KIT INFORMATION: 1. Test kit used for this sample: Yes, No
2. Results: PCB not present; PCB present, less than 50 ppm; PCB present, greater than 50 ppm; _____ 100 % PCB present. SPILL-FYTER CHEMICAL CLASSIFIER STRIPS: Results: Chloring Risk

+ for to ton Driger Tube **COMMENTS**

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 1 9150

Samplers: Samplers: Samplers: Samplers: Samplers: Sample Location: Sample	REAC Task Leader: Rae EPA WAM: Gores	Chain of Custody No.: 3627 REAC Task Leader: Ray Lean; EPA WAM: Gres Forcy Work Assignment No.: 33474001000801			
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	SOIL TYPE SURFACE W/ rock clay color gravel muck odor sand loam flow silt peat direction color	width depth velocity cm/s	BOTTOM rock silt rubble clay gravel organic shell other sand		
SAMPLE TYPE DEV surface water effluent kemi groundwater sludge trowe potable water léachate buck sediment waste auge soil otherekma	nerer ponar colo other odor et temp	sample depth	- ' 		
ANALYSES TO BE PERFORMED ORGANICS 1. halogenated & aromatic volatiles B. volatiles C trihalomethanes D pesticides/PCB E. PCB F. base neutral/acid extractables G pesticides, drinking water H. herbicides, drinking water I other INORGANICS A. metals, priority pollutant B metals, TAL C metals scan (ICP) D metals, other	OTHER ANALYSES A. total cynanide B. total phenol C. petroleum hydrocarbo D. pH E. alkalinity F. hardness G. total dissolved solids H. total suspended solids I. sulfate J. TOC K. grain size L. percent moisture M. other	plastic bag plastic bucket other	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other		
RCRA A TCLP B. ignitability C. corrosivitypH D. reactivity E other					

Drun Sunple REAL#9
FLA DEP # 9
5: vry Om 1714/

FORM #1

REAC# 10 100 PLADEP# Photo 20 17E	
Om is don't in the side. writing on top of Drum "407# MCC 22-23	
Dong is full, Sould Blue Worght Anitals	
- Ream of January Strik	
	

2 2 3 3 3 2

Samplers: Getty & Morganti

Date: 10 Jan 95

Site Name: Pier Drum

Work Assignment Number: 3347-04-01-0098 REAC Task Leader: Ray Lewis

Container Number/Sample Number: REAC#10

SITE INFORMATION:

Terrain, drainage description: Flat, roul woods and feilds with sandy soil.
 Weather conditions (from observation): Mid to high 60's °F and sunny.

MET station on site: X No, Yes

SAMPLE INFORMATION

	SAMPLE INFORMATION:
1.	Container type: X Drum, Tank, Other:
2.	Container dimensions: Shape: 17E in overpak-17H Approximate size: 55 gal in 85 gal
3.	Label present: X No, Yes:
4.	Spill or Leak present: X No, Yes, Dimensions:
	Description: liquid, X solid (powder or X crystals), sludge
6.	Color: , Odor: Yes , Vapors: Yes
	FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING
	Flash No Flash
	VI OF LOUROUS COLUMNOS

pH OF AQUEOUS SOLUTION 1. Using 0-14 pH paper, check pH of water/sample solution:

WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATER REACTIVE!] Results: _____ total, ____ partial, ____ no solubility.

•	
WATER REACTIVITY:	
1. Add small amount of sample to water: bubbles, color change to	_,
odor or vapor formation, heat, No Change.	_
SPECIFIC GRAVITY TEST (compared to water):	
1. Add small amount of sample to water: X Sinks, Floats.	
2. If liquid comple sinks coroon for obloginated compounds. If liquid comple floats and appears to	_

2. If the sid seconds, sinks			If the said seemeds	floats and ammass	es to be silv
2. If liquid sample sinks,	, screen for chlorina	tea compounds.	ii iiquid sampie	moats and appear	is to be only,
screen for PCB's (Chlo	or n Oil kit).				

CULOD I OU TEST AIT INFORMATION.

	CHECK NOL TEST KIT INTOKMATION.
1. Test kit used	for this sample: Yes No
2. Results:	PCB not present; PCB present, less than 50 ppm;
	PCB present, greater than 50 ppm; 100 % PCB present.

SPILL-FYTER CHEMICAL CLASSIFIER STRIPS:

Results:

+ for Styn Drig - Tube COMMENTS

00610

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022 2 1 11113

Samplers:	men	REA		BOTTOM rock silt rubble clay gravel organic shell other sand
SAMPLE TYPE DEVICE surface water effluent kemmer groundwater sludge trowel potable water leachate bucket sediment waste auger soil other ekman		SAMPLE INFO color_ odor_ temp DO_ cond	pH	WEATHER PARAMETERS ambient temp_ barometric pressure_ relative humidity_ weather conditions
ANALYSES TO BE PERFORMED	 		SAMPLE PREPAR	
ORGANICS 1. halogenated & aromatic volatiles 2. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water I other INORGANICS A. metals, priority pollutant B. metals, TAL C. metals scan (ICP) D. metals, other	A. total B. total C. petr D. pH E. alka F. hard G. tota H. tota I. sulfa J. TOO K. grain L. perd	Iness I dissolved solids I suspended solids te	container glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
RCRA A. TCLP B. ignitability C. corrosivitypH D. reactivity E. other				

Drum Sample REACHTIQ FLADER #10 17E

FORM #1

2 2 7734

		30
0.	FAC # 11	
- 1/	EXC#11 FLAPEP# E	Photo 21
17	ε	
		I Point is porting
	Don is full h	,1 Solid Rod +
	- Clay	
	g (F), i i i i i i i i i i i i i i i i i i	
<u></u>		
* - · ·		a commence of the second of th
	··	

10 Jan 95 Samplers: Getty & Morganti Date: Work Assignment Number: 3347-04-01-0098 Site Name: Pier Drum REAC Task Leader: Ray Lewis Container Number/Sample Number: REAC#11 SITE INFORMATION: 1. Terrain, drainage description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's °F and sunny. MET station on site: X No, Yes SAMPLE INFORMATION: 1. Container type: X Drum, Tank, Other: 2. Container dimensions: Shape: 17E in overpak 17H Approximate size: 55 gal in 85 gal 3. Label present: X No. Yes: 4. Spill or Leak present: X No, Yes, Dimensions: X solid (powder or X crystals), sludge 5. Description: liquid, 6. Color: , Odor: Yes , Vapors: Yes FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING **IGNITABILITY** No Flash pH OF AQUEOUS SOLUTION 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATERREACTIVE!] Results: _____ total, ____ partial, ____ no solubility. WATER REACTIVITY: 1. Add small amount of sample to water: _____ bubbles, ____ color change to _____ odor or vapor formation, _____ heat, ____ No Change. SPECIFIC GRAVITY TEST (compared to water): 1. Add small amount of sample to water: X Sinks, ____ Floats. 2. If liquid sample sinks, screen for chlorinated compounds. If liquid sample floats and appears to be oily, screen for PCB's (Chlor n Oil kit). CHLOR N OIL TEST KIT INFORMATION: 1. Test kit used for this sample: PCB not present; PCB present, less than 50 ppm;
PCB present, preater than 50 ppm; 100 % PCB present. 2. Results: SPILL-FYTER CHEMICAL CLASSIFIER STRIPS:

COMMENTS

+ for tester Driger Tute

Results:

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022



	Samplers:	er Pr	expute m obe Pet		Chain of Custody No.: 3 REAC Task Leader: 4 EPA WAM: 5 Work Assignment No.: 3	C 27 / Lews Evel Y7-4001000880V
SITE DESC landfill industrial commercial residential hedgerows	old field upland palu wooded lowland rive		SOIL TYPE rock clay gravel muck sand loam silt peat color	SURFACE WATER color odor flow direction		BOTTOM rock silt rubble clay gravel organic shell other
SAMPLE TY surface water groundwater potable water sediment soil	er effluent sludge	DEVICE kemmere trowel bucket auger ekman	r ponar other	SAMPLE color odor temp DO cond	ORPsalinitysample_depth	WEATHER PARAMETERS ambient temp_ barometric pressure_ relative humidity_ weather conditions
ORGANICS A halogenat B volatiles C trihalomet D pesticides E PCB F base neut G pesticides H herbicides	thanes s:/PCB tral/acid extractables s: drinking water s: drinking water		A. tota B. tota C pet D pH E. alka F hard G tota H tota J TOO	dness al dissolved solids al suspended solids ate	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
B. metals, TAC metals scoop metals, of metal	an (ICP) ther		L. per	cent moisture	ambient	

COMMENTS

Drun Sample REAC # 11
PLA DEP # 11

FORM#1 17E

	REHC # 12
	17 FLADER # Photo 22
	D. : : 6./ =/ 1.
	Don is in book 5hope.
	The Top is Coushed.
	(Photo 23)
	$0.1 \cdot 1 \leq 1 \leq 1$
	And _ Somphy Som 25
	<i>+</i> _/ <i>q</i>
144	
-	
٠	
••	
	<u> </u>
>	

			37	1 2 1
REAC # 13	7 	0/.1. >=		
13 FLAD 17E	ε P	PAOTO_23_		
Don is i	in for 5h	pr		
3/4 Full Onither	w/ Soul a	green ? way	jat	
•				
	. .	and the second second second second second		
• • • • • • • • • • • • • • • • • • •				
·• ·				
		- a 		
			· · · · · · · · · · · · · · · · · · ·	
and the second of the second of the second of				

DRUM/TANK SAMPLING DATA SHEET

10 Jan 95

Date:

Samplers: Getty & Morganti Work Assignment Number: 3347-04-01-0098 Site Name: Pier Drum Container Number/Sample Number: REAC#13 REAC Task Leader: Ray Lewis SITE INFORMATION: 1. Terrain, drainage description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's F and sunny. MET station on site: X No, Yes SAMPLE INFORMATION: 1. Container type: Tank, Other: X Drum, 2. Container dimensions: Shape: 17E in overpak 17H Approximate size: 55 gal in 85 gal 3. Label present: X No. Yes: 4. Spill or Leak present: X No, Yes, Dimensions: 5. Description: liquid. X solid (powder or X crystals), sludge 6. Color: , Odor: Yes , Vapors: Yes FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING **IGNITABILITY** No Flash Flash pH OF AQUEOUS SOLUTION 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATERREACTIVE!] Results: _____ total, ____ partial, ____ no solubility. WATER REACTIVITY: 1. Add small amour of sample to water: _____ bubbles, ____ color change to _____ no Change. SPECIFIC GRAVITY TEST (compared to water): 1. Add small amoun' of sample to water: Sinks, ____ Floats. 2. If liquid sample onks, screen for chlorinated compounds. If liquid sample floats and appears to be oily, screen for PCB's (Chlor n Oil kit). CHLOR NON TEST KIT INFORMATION:

1. Test kit used for this sample: Yes, No

2. Results: PCB not present; PCB present, less than 50 ppm; PCB present.

PCB present, greater than 50 ppm; 100 % PCB present. SPILL-FYTER CHEMICAL CLASSIFIER STRIPS: Results:

COMMENTS

· for Styren Dorger Tube

00613

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 2 3.03

Samplers:	puz .	REAC Task Leader: Repair Repai	rel/
SAMPLE TYPE DEVICI surface water effluent kemme groundwater studge trowel potable water leachate bucket sediment waste auger soil other ekman		EINFORMATION pHORPsalinitysample_depth	barometric pressure
ORGANICS A. halogenated & aromatic volatiles B. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water I. other INORGANICS A. metals, priority pollutant B. metals, TAL C. metals, other D. metals, other	OTHER ANALYSES A. total cynanide B. total phenol C. petroleum hydrocarbons D. pH E. alkalinity F. hardness G. total dissolved solids H. total suspended solids I. sulfate J. TOC K. grain size L. percent moisture M. other	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
RCRA A TCLP B ignitability C corrosivitypH D reactivity E other COMMENTS Drum Sumply FLA DEP # 1-	REAC# 13		

PERCH 14 14 FLADER Phuk 24 17E DM 13 Out of Ships. Full w/ Gran 4 Weight Solid Christles.
DM 13 out of Ships
Full w/ Grands - Weight Sold Corales -
Sold Cristles.
50 W 073 NO

DRUM/TANK SAMPLING DATA SHEET

9162 10 Jan 95 Samplers: Getty & Morganti Date: Work Assignment Number: 3347-04-01-0098 Site Name: Pier Drum REAC Task Leader: Ray Lewis Container Number/Sample Number: REAC#14 SITE INFORMATION: 1. Terrain, drainage description: Flat, roul woods and feilds with sandy soil. 2. Weather conditions (from observation): Mid to high 60's °F and sunny. MET station on site: X No. Yes SAMPLE INFORMATION: 1. Container type: X Drum, Tank, Other: 2. Container dimensions: Shape: 17E in overpak 17H Approximate size: 55 gal in 85 gal 3. Label present: X No. Yes: Yes, Dimensions: 4. Spill or Leak present: X No, liquid, X solid (5. Description: powder or X crystals), sludge 6. Color: , Odor: Yes , Vapors: Yes FIELD TEST DATA SHEET FOR DRUM/TANK SAMPLING **IGNITABILITY** No Flash Flash pH OF AQUEOUS SOLUTION 1. Using 0-14 pH paper, check pH of water/sample solution: WATER SOLUBILITY TEST: 1. Add approximately one part sample to five parts water. You may need to stir and heat gently. [DO NOT HEAT IF WATERREACTIVE!] Results: _____ total, ____ partial, ____ no solubility. WATER REACTIVITY: 1. Add small amount of sample to water: _____bubbles, ____ color change to _____, odor or vapor formation, _____ heat, ____ No Change. SPECIFIC GRAVITY TEST (compared to water): 1. Add small amount of sample to water: ____ Sinks, ___ Floats. 2. If liquid sample sinks, screen for chlorinated compounds. If liquid sample floats and appears to be oily, screen for PCB's (Chlor n Oil kit). CHLOR N OK TEST KIT INFORMATION: 1. Test kit used for this sample: PCB present, greater than 50 ppm; _____ 100 % PCB present. 2. Results:

COMMENTS

SPILL-FYTER CHEMICAL CLASSIFIER STRIPS:

+ for Styres Driger Tobe

Results:

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022 2 1 1165

	ipie Lucation	Singo	Samplers: Craffy Mcgashi Date Mark: Rei Driv Time: 1300 Sample Location: Sansata				
SITE DESCRIPTION landfill old field industrial wooded commercial farmland residential gully hedgerows floodplain	upland palusti lowland riverii lacustrine	rine ne	gravel sand	clay muck loam peat	SURFACE WATER color odor flow direction	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other sand
sediment was	ge hate	DEVICE kemmerer trowel bucket auger ekman		ar er	SAMPLE color odor temp DO cond	ORPsalinitysample_depth	WEATHER PARAMETERS ambient temp barometric pressure relative humidity weather conditions
ORGANICS A. halogenated & aroma J. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extr G. pesticides, drinking w H. herbicides, drinking w I. other	tic volatiles actables ater ater			A. total community B. total pi C. petrole D. pH E. alkalin F. hardne G. total di	henol eum hydrocarbons ity	SAMPLE PREPA CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
NORGANICS A metals, priority polluta B. metals, TAL C. metals scan (ICP) D. metals, other	nt			K. grain s L. percen	ize t moisture	dry ice ambient	
RCRA A TCLP B ignitability C corrosivity D reactivity E other	 						

REAC #15	
15 PLANEP # PM	ut 25
Dm 13 doubed on to	hy siche:
Rock Solid to the	top, could not
5 impher	
······································	
· · · · · · · · · · · · · · · · · · ·	

REAC#16
16 FLADEP # Photo 26
52breve 17H W/17E in suche
NA Somples
NA Somples Rock Sold to the Top
•
<u> </u>

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 1 3165

Date: 30 Jan 95 Site Name: New Sample Location:	Down 30750fr FLA	REA	n of Custody No.: C Task Leader:	115 11 10088 m 47-410010088
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	rock clay color _ gravel muck odor _ sand loam flow	ACE WATER	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other sand
SAMPLE TYPE DEVICE surface water effluent kemm groundwater sludge trowel potable water leachate bucke sediment waste auger soil other ekmar	erer ponar other	SAMPLE INFO color odor temp_ DO cond	pH	barometric pressure relative humidity weather conditions
ANALYSES TO BE PERFORMED ORGANICS A. halogenated & aromatic volatiles 3. volatiles C. trihalomethanes D. pesticides/PCB E. PCB F. base neutral/acid extractables G. pesticides, drinking water H. herbicides, drinking water I. other	OTHER ANALYS A. total cynanide B total phenol C. petroleum hydi D pH E. alkalinity F hardness G total dissolved H. total suspende	rocarbons solids	SAMPLE PREPAR CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
INORGANICS A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals other VOA RCRA A TCLP	J. TOC K. grain size L. percent moistu M. other		wet ice dry ice ambient	
B ignitability C corrosivity D. reactivity E other				

FORM #1

REACH 8
REACH 9
REACH 4
REACH 4

REAC, EDISON, NJ (908) 321-4200 **EPA CONTRACT 68-C4-0022**

Samplers: Getty Date: 30-12495 Site Name: Rev Drug Time 100 Sample Location: PEHC	Chain of Custody No.:	us s U
SITE DESCRIPTION SOIL TYPE SURFACE WATER and fill old field upland palustrine rock clay color rock clay color gravel muck odor sand loam flow flow silt peat direction color floodplain color color color floodplain flood	width depth velocity cm/s	BOTTOM rock silt rubble clay gravel organic shell other
SAMPLE TYPE DEVICE SAMPLE surface water effluent kemmerer ponar color groundwater sludge trowel other odor_ potable water leachate bucket temp_ sediment waste auger DO_ soil other ekman cond ANALYSES TO BE PERFORMED	ORPsalinitysample_depthtide_stage	barometric pressure relative humidity weather conditions
ORGANICS A balogenated & aromatic volatiles Volatiles C. trihaiomethanes D pesticides/PCB E. PCB C. petroleum hydrocarbons D pH E. alkalinity F. hardness G pesticides, drinking water H herbicides, drinking water H. other INORGANICS A metals, priority pollutant B metals, TAL OTHER ANALYSES A total cynanide B. total phenol C. petroleum hydrocarbons D pH E. alkalinity F. hardness G total dissolved solids I. sulfate J. TOC K. grain size L. percent moisture M other M other	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient A S III	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other

Sample was Split between 2 Contract Lab REACCES

REAC, EDISON, NJ (908) 321-4200 EPA CONTRACT 68-C4-0022

2 . 0163

Samplers:	Orun Et C	RE	A WAM: Pur	66 LUIS LEU BY 140010008
SITE DESCRIPTION landfill old field upland palustrine industrial wooded lowland riverine commercial farmland lacustrine residential gully hedgerows floodplain	SOIL TYPE rock clay gravel muck sand loam silt peat color	SURFACE WATER color odor flow direction	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other sand
SAMPLE TYPE DEVICE surface water effluent kemm groundwater sludge trowel potable water leachate bucket sediment waste auger soil other ekmar	erer ponar other	SAMPLE INFO color odor temp DO cond	pH	
ORGANICS A halogenated & aromatic volatiles 3 volatiles C trihalomethanes D pesticides/PCB E PCB F base neutral/acid extractables G pesticides drinking water H herbicides, drinking water I other	A. total c B. total p C. petrole D. pH E. alkalin F. hardne G. total d	henol eum hydrocarbons ity	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other
INORGANICS A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals other RCRA A TOLD B ignitability Metals	K. grain s L. percer	size nt moisture	wet ice dry ice ambient	

COMMENTS

Sample from Down REAL# 5

REAC, EDISON, NJ (908) 321-4200 **EPA CONTRACT 68-C4-0022**

2 1 0.39

Samplers: 7. Date 301 15 Site Name: 1 Time: 1010 Sample Location:	111	RE RE EP	Chain of Custody No.: REAC Task Leader: Leal i EPA WAM: Powell Work Assignment No.: 334740000088		
SITE DESCRIPTION landfill old field upland palus industrial wooded lowland river commercial farmland lacustrine residential gully hedgerows floodplain	•	SURFACE WATER color odor flow direction	STREAM width depth velocity cm/s pools% riffles%	BOTTOM rock silt rubble clay gravel organic shell other sand	
SAMPLE TYPE surface water effluent groundwater sludge potable water leachate sediment waste soil other	DEVICE kemmerer ponar trowel other bucket auger ekman	SAMPLE INF color_ odor_ temp DO_ cond	pHORPsalinitysample_depth	barometric pressure	
ORGANICS A. halogenated & aromatic volatiles 3. volatiles C. trihalomethanes D pesticides/PCB E PCB F base neutral/acid extractables G pesticides drinking water H herbicides drinking water I other INORGANICS A metals, priority pollutant B metals, TAL C metals scan (ICP) D metals, other RCRA BUA TCLP PCO/PAT B ignitability C corrosivitypH D reactivity E other	A. total B. total C. petro D. pH E. alkal F. hardi G. total H. total J. TOC K. grain L. perce	oleum hydrocarbons linity ness dissolved solids suspended solids	CONTAINER glass jar plastic jar acetate core plastic bag plastic bucket other STORAGE wet ice dry ice ambient	PRESERVATIVES HNO, NaOH Zn Acetate HCI Na,SO, other	

Composit Sample from Drim Samples REAC#3
REAC#10
PENE#13
PO10 # 14

FORM#1

2080.r99 8 of 99.9 11/12/92

TOXICITY CHARACTERISTIC LEACHING PROCEDURE (TCLP)

TCLP Sample Data Sheet

		EXTRACTION	CTION	FLUID DETERMINATION	ATION		TCLP	TCLP EXTRACTION	7	
SAMPLE # AND DESCRIPTION	% WET SOLIDS	FILTRATE pH (%solid <100)	INITIAL pH FINAL pH (solids)	FINAL pH (solids)	EXTRACTION FLUID	FLUID VOLUME	SAMPLE MASS	START TIME	END TIME	FINA pH
A 00642	83		(0.3	h'1	#	11	150/11	1520	0925 4.99	4.9
A 00643	83		6.2	1.6	1#	1	50 /m/		_	5.0
A 00644	83		6.3	1.3	1#	2r	100(120)			4.8
A 60 42B	/ . \				#	11	50 (4)			4.93
438					#	_ 1F_	50 (41)			5.0
d 44 B				/	1#	26	100(120)	,		4,9
420			X		#	11	(19) 05			5.0
784					#	ור	(19)05			5.0
244					#	7K	(021) (201	>	→	16'h
			•				\ -			
					,	×				
	J 4 1 . •								2	

Technician:
Date: //CC-C-C-R:

Site Name: Pier N. WA #:
Task Leader: Run L. WAM:

0.710

CHAIN OF CUSTODY RECORD/LAB WORK REQUEST

Roy F.: ston, Inc. REAC, Edison, N.J. EPA Contract 68-03-3482

No: 10027

				,			,		,					,		•		 		12 2		9121
												-		_	·			 	, , ,	ONLY	Time	
		_					-												7	NG USE	Date	
TCL) Pest			X			У,			X									1	<i>I</i>	FOR SUBCONTRACTING USE ONLY FROM CHAIN OF CUSTODY #	Received By	
ENS		X			X			X										1		FOR SUBCO FROM CHAIR CUSTODY #	Date R	
11-4-15	X			χť			7														Relinquished By	
Container/ Preservative	L 12 Lu / 112 118 1	_ `4	11. 1 mbe 1916	1. Tale / 11 1 V.	1 Posts 14 c	•	L 156,79 ar	1. 1 milion / 1/10	1 Purbar/11 c					,			,			(Is)	Items/Reason R	
# or Bottles	۲,				-				P -				_							Instructions: E F = [15,175]	Time	08. P
Date Collected	1/18/15	-							5					/						oecial Instruct		11 21/8/1
Matrix	\^\								احتوا											Soil Special Ir water (1952-7 19, E Oil Air	Received By	Marker (M)
Sampling Location	64.2			< μ.		-	1)													თ≯О∢	Date	19 Car Carlo
Samplin	ACCOUR	-	1.2	100 64	-	_	1-200 V		-		1									Potable Water Groundwater Surface Water Sludge	Relinquished By	7000
Sample No.	50500	1300 S.C.	5.5000	9 85004	9750	1 .	1.5000	B.T. (1977)	0.522											Sediment PW - Drum Solids GW - Drum Liquids SW - Other	_	Allel cirpland
REAC #	V				_	.	<	<u> </u>	! O	,				Vandage			 			Matrix: SD - Sedim SS - Drum CL - Drum C - Other	Items/Reason	11.1

EPA Contract 68-C4-0022 . REAC, Edison, NJ (908) 321-4200

CHAIN OF CUSTODY RECORD

-

33174041001801801 Project Number. Project Name:_

03627 ë

SHEET NO. ZOF 1

Analyses Requested 6. Container/Preservative # of Bottles Special Instructions: Date Collected Sample Identification Matrix 31/05 \$ 0 € *D* \$ d á 80 Potable Water Groundwater Surface Water Sludge Sampling Location JM1/1 0,413 041/1 11115 £ %. %. %. Sample No. 600 2005 Sediment Drum Solids Drum Liquids Other REAC# SD д×

FOR SUBCONTRACTING USE ONLY FROM CHAIN OF **CUSTODY#**

·No: 03663 FOR SUBCONTRACTING USE ONLY SHEET NO._OF_ 8 Pest/Reb X St. Analyses Requested Received By FROM CHAIN OF CUSTODY # X Metho Relinquished By 03890100000000000850 Container/Preservative 74.6 Items/Reason TANK # of Bottles Special Instructions: Time 4 4 Project Number.
RFWContact: 4 M Data Collected Sample Identification Received By Matrix Potable Water Groundwater Surface Water Sludge Sampling Location Relinguished By EPA Contract 68-C4-0022 . SW. SW. 1150041 500041 Sample No. 4 कि के हैं न F00641 POOLY Sediment Drum Solids Drum Liquids Other Nema/Reason REAC# ORM #4 Metric: SD-

CHAIN OF CUSTODY RECORD

Project Name:

REAC, Edison, NJ (908) 321-4200 03664 FOR SUBCONTRACTING USE ONLY E E SHEET NO COF **\$** .. Z **Analyses Requested** Received By FROM CHAIN OF CUSTODY# TCLP VOAS Relinquished By 03347040001009861 Phone: Container/Preservative Items/Reason 200 # of Bottles Special Instructions: Time Date Collected 56/11/ Soil Water Oil Air Sample Identification Received By Martrix Potable Water Groundwater Surface Water Sludge Sampling Location Relinquished By **EPA Contract 68-C4-0022** ₽ \$ \$ \$ \$ Sample No. 300002 800613 Sediment Drum Solids Drum Liquids Other thems/Reason A Mandys. 3 REAC# FORM #4

小道 人名英格兰人

H

CHAIN OF CUSTODY RECORD

Project Name:__

REAC, Edison, NJ

(908) 321-4200

REAC, cdison, NJ (908) 321-4200 EPA Contract 68-C4-0022

CHAIN OF CUSTODY RECORD

The section of the second section of

a contract the second of the second

1334704000100980 Phone: ur Irum Project Name:

03870 ë

SHEET NO. OF

		Sample Identification	entifica	tion			Analy	Analyses Requested	SHEET NOOF	7.—GF——
REAC #	Sample No.	Sampling Location	Matrix	Date Collected	semog jo #	Container/Preservative	Mitale	BNA	Pesters	2, 401
	4600635		ω	56/11/1	7	HNO, 14°C	×		,	
	1000635		1	1,	c	7.00		У		
	FF00635			1	7	,			×	
_	642 00635				3					×
	4320659				2	1100 /40C	×			
	D00639				C	7.4.		×		
	FF00639				2				×	
	CHI 00639				8					×
-	01900gv				2	4°C/HNU?	X			*
) _	96 90 9 A3				2	Jah		×		
	5F 00 640				7	1			X	
t -	9/1900 IHS		1	1	3					×
1									,	
		/								
		1								
 -	-	•								
	/									
	e									-
×	Sediment	PW - Potable Water	v.		Special Instructions:					
08-10	sbi sbi		န်ဝံ	Water			FOR S	UBCONTR	FOR SUBCONTRACTING USE ONLY	SE ONLY
			¥	Air			1000	EDOM CUAIN OF		

FOR SUBCONTRACTING USE ONLY FROM CHAIN OF CUSTODY #

ftems/Reason	Refinquished By	Date	Received By	Date	Time	Items/Reason	Relinquished By	Date	Received By	Date	Time
Juffer to	1	11.11.5									
	,							,			
FORM #4											8/94

03671 FOR SUBCONTRACTING USE ONLY SHEET NO._OF_ **YoA'S** Q SE AH/RBS .; **2** Analyses Requested Received By FROM CHAIN OF CUSTODY # Date Relinquished By 8600 100 040 Phone: Container/Preservative) (Items/Reason **₩** # of Bottles Special Instructions Time UM Project Number: Project Name:_ 13/95 RFW Contact: Date Date Collected Sample Identification Received By Matrix Potable Water Surface Water Surface Water Surface Sampling Location Date Relinquished By EPA Contract 68-C4-0022 1-AI 00636 ABOOK 38 MT 20638 20000 4800636 F00636 D00638 F10638 Sample No. · REAC, Edison, NJ Sediment Drum Solids Drum Liquids Other (908) 321-4200 Items/Resson 410 FORM #4 13. <u>\$</u>1, *

CHAIN OF CUSTODY RECORD

03672 FOR SUBCONTRACTING USE ONLY SHEET NO._OF_ 246 Motole PraiRB's BNA'S X Received By AC Analyses Requested FROM CHAIN OF CUSTODY # X \$ Relinquished By Phone: Brian Holdenis 03900 100 040 THES Container/Preservative Items/Resson HNOZ HAIDZ 000 Soil Nation Far A00643 oil A00644 Air A00642 # of Bottles Special Instructions Time JAGAN AM Project Number. RFW Contact: Date Collected Sample Identification Received By Matrix 3 S Potable Water Groundwater Surface Water Sludge Sampling Location 1/4-1/5 Relinquished By EPA Contract 68-C4-0022 100642 THE 00642 HT 00657 #200639 300637 EF 00637 A 200643 Vraugh 4 Sample No. 10643 80063 D00634 1600657 600631 100644 Sediment Drum Solids Drum Liquids Other Hems/Reason REAC# FORM #4

CHAIN OF CUSTODY RECORD

Project Name:__

REAC, Edison, NJ

(908) 321-4200

REAC, Edison, NJ (908) 321-4200 EPA Contract 68-C4-0022 *

CHAIN OF CUSTODY RECORD

Project Number.
RFW Contact: / Project Name:__

No: 00064

Phone:

SHEET NO. LOF _

J		Π	Γ	١,	Γ	T			Γ	T		Π	Г				1	-	Τ
						1												SE ONL	
pa		\						\										CTING US	
Analyses Requested			\															FOR SUBCONTRACTING USE ONLY	
Analyse	KON	7	7	7					/	\								FOR SU	
	Container/Preservative	Yell I'm Pl	/	//															
	# of Bottles	,	,	,													Special Instructions:		
tion	Data Collected	1 Jn. 15	16 10.13	11/ 1							,						ŀ	Water	5 ₹
entifica	Metrix	70	70	70													,	5 \$ ċ	►
Sample Identification	Sampling Location	11.11	54. HZ	1,# 10											<u> </u>		l	Groundwater Surface Water	
	Sample No. Sa	┤	4.0413 V															Drum Solids GW-	
500510	REAC #	1569	p 662	300									_			\	Matrix		

FOR SUBCONTRACTING USE ONLY **FROM CHAIN OF CUSTODY#**

	Nema/Resson	Relinquished By	Date	Received By	Date	Time	Time Items/Reason	Relinquished By Date	Date	Received By	Date	Time	_
T.	RILLIE	MIN 1	1	1 Lewan	130/95	1:05	4:05 3/ VOA						
*-	77177				/			•					
<u> </u>													_
<u>. </u>													_
													_
J2	Sobre 44											3	-

00065 SHEET NO. LOF L FOR SUBCONTRACTING USE ONLY Melel Ö 130/2 1 Analyses Requested Received By FROM CHAIN OF PCB/Lit CUSTODY # Phone: 4 13/1/17 de Ž Relinquished By Container/Preservative Analysis Items/Reason 1 # of Bottles Special Instructions: 20 Project Number.
RFW Contact: 136/95 Project Name:_ Date Collected ř Solf Vater ≥ ⊆ Sample Identification Received By Metrtx 70 Potable Water Groundwater Surface Water Sludge Sampling Location DENCE CI サンプラ Relinquished By EPA Contract 68-C4-0022 Sample No. REAC, Edison, NJ (908) 321-4200 121 600 1600 Logg! Sediment Drum Solids Drum Liquids Other · REAC 1

PORMI

CHAIN OF CUSTODY RECORD

99000 SHEET NO.L OF L FOR SUBCONTRACTING USE ONLY Ö Received By Analyses Requested FROM CHAIN OF **CUSTODY#** 21416 Relinquished By Phone: 12. Containen Preservative, 18 18 1 C ttems/Reason Special Instructions: # of Bottles Ē Project Number: RFW Contact: Project Name: **Data Collected** 1 Soil Water Oil Air Received By Sample Identification 37.1 DY Matrix ÷ 0 ₹ 0.+ $\stackrel{\sim}{\sim}$ Potable Water Groundwater Surface Water Sludge Sampling Location 8 950 67 # 75 Relinquished By **EPA Contract 68-C4-0022** SW-SW-SL-SIL-Sample No. Sediment Drum Solids Drum Liquids Other (908) 321-4200 Nems/Resson 013095 REAC# ORM #4 و SD.

REAC, Edison, NJ

1

CHAIN OF CUSTODY RECORD

Appendix B

APPENDIX B

FIELD SCREENING RESULTS
Pier Drum Site
Bradenton, FL
January 1995

Appendix Contents:

Laboratory HazCat Data

DRUM INVENTORY LOG

orum Id. : 1

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Гime

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 75%

Drum Cond.: POOR

Overpack Size:

Debris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	SOLID	BROWN		
2	(Middle)				
3	(Bottom)				

Manufacturer:

Chemical :

Guerator:

Misc. : The bung is moist and may be leeking

LABORATORY HAZCAT DATA

2 : 0:04

Drum ID : 1

Li 2r : (Top)

Color : BROWN

Physical State: SOLID

Clarity :

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	
WATER REACTIVITY	N	
RADIOACTIVITY	x	
WATER SOLUBILITY	Y	Partial
HEX/DCM.SOLUBILITY	N	
рн	5	pH is of the water used in solubility test
OXIDIZER	х	
PEROXIDE	x	
CYANIDE	x	
LLFIDE	x	
CHLORIDE	N	
FLAMMABLE	Y	sample flashed after 1 min of heating
РСВ	N	

DRUM INVENTORY LOG

2 2 0:05

Drim Id. : 2

Location: pit

ate : 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

rum Top : BUNG

% Full : 100%

Drum Cond.: FAIR

Overpack Size:

Jebris/PPE:

No. of Layers: 2

-	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	LIQUID	BROWN	CLOUDY	50%
2	(Middle)				
13	(Bottom)	SLUDGE	BLUE	CLOUDY	50%

'Manufacturer:

C._mical :

Generator :

Misc. : Markings: D10N Polyester Resin

LABORATORY HAZCAT DATA

Drum ID : 2

I er : (Top)

Physical State: LIQUID

Color : BROWN

Clarity : CLOUDY

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	
WATER REACTIVITY	N	
RADIOACTIVITY	x	
WATER SOLUBILITY	Y	Partial
HEX/DCM SOLUBILITY	х	
рн	7	
OXIDIZER	x	
PEROXIDE	х	
CYANIDE	x	
, JLFIDE	x	
CHLORIDE	x	
FLAMMABLE	x	Chlorine wine test produced a green flame
PCB	Y	Quanti Chlor Test yielded from 844ppm to 2400 ppm

LABORATORY HAZCAT DATA

orum ID : 2

L er : (Bottom)

Physical State: SLUDGE

Color : BLUE

Clarity : CLOUDY

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	N	
WATER REACTIVITY	N	
RADIOACTIVITY	N	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	N	
рН	N	
OXIDIZER	N	
PEROXIDE	N	
CYANIDE	N	
JLFIDE	N	
CHLORIDE	N	
FLAMMABLE	N	
PCB	N	

DRUM INVENTORY LOG

2 1 0133

Drum Id. : 3

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 75%

Drum Cond.: POOR

Overpack Size:

Debris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	SOLID	AMBER	CLOUDY	100%
2	(Middle)				
3	(Bottom)				

Manufacturer:

C. _mical :

Generator :

Misc.

: Plastic like, very hard to sample

ADDITIONAL DRUM INFORMATION

Drum Class. :

Bulk Number :

Disposal Opt.:

Disposal Co. :

Manifest No. :

LABORATORY HAZCAT DATA

rum ID : 3

'. er : (Top)

Physical State: SOLID

Color : AMBER

Clarity : CLOUDY

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	Positive for Styrene Drager Tube
WATER REACTIVITY	N	
RADIOACTIVITY	x	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	x	
рН	7	
OXIDIZER	x	
PEROXIDE	x	
CYANIDE	x	
JLFIDE	х	
CHLORIDE	x	
FLAMMABLE	х	
PCB	х	

DRUM INVENTORY LOG

Drum Id. : 4

Location: pit

Date : 01/10/95

Sampler : S.Getty\M.Morganti

Time :

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17H

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 25%

Drum Cond.: POOR

Overpack Size: 55-GAL

Debris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	LIQUID	COLORLESS	CLEAR	
2	(Middle)				
3	(Bottom)				

Manufacturer:

C. mical :

Generator :

Misc.

LABORATORY HAZCAT DATA 2 1 3491

rum ID : 4

T er : (Top)

Physical State: LIQUID

Color : COLORLESS

Clarity : CLEAR

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	х	
WATER REACTIVITY	N	
RADIOACTIVITY	x	
WATER SOLUBILITY	Y	
HEX/DCM SOLUBILITY	x	
рн	7	
OXIDIZER	х	
PEROXIDE	х	
CYANIDE	х	
JLFIDE	х	
CHLORIDE	x	
FLAMMABLE	N	
РСВ	х	

Drim Id. : 5

Location: pit

Date : 01/10/95

Sampler : S.Getty\M.Morganti

Time :

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17H

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 100%

Drum Cond.: POOR

Overpack Size: 85-GAL

Debris/PPE:

No. of Layers: 1

LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1 (Top)	GEL	COLORLESS	CLEAR	100%
2 (Middle)				
3 (Bottom)				

Manufacturer:

Cilmical

Generator :

Misc.

2 2 3193

)rum ID : 5

L_ar : (Top)

Physical State: GEL

Color : COLORLESS

Clarity : CLEAR

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	
WATER REACTIVITY	N	
RADIOACTIVITY	x	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	x	
рН	7N/A	
OXIDIZER	х	
PEROXIDE	x	
CYANIDE	x	
LFIDE	x	
CHLORIDE	x	
FLAMMABLE	N	
PCB	х	

Drum Id. : 6

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 75%

Drum Cond.: FAIR

Overpack Size:

Debris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)				
2	(Middle)				
3	(Bottom)				

Manufacturer:

C. .nical

Generator :

Misc.

: Contents appear to have polymerized. No sample.

2 1 0:25

Drum Id. : 7

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time :

Witness : S.Getty\M.Morganti

Drum Type : BLUE AND WHITE DRUM

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 75%

Drum Cond.: POOR

Overpack Size:

Debris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	SOLID	PURPLE	CLOUDY	100%
2	(Middle)				
3	(Bottom)				

Manufacturer:

C. _mical :

Generator :

Misc. : Very hard to sample

Drum ID : 7

La r : (Top)

Physical State: SOLID

Color : PURPLE

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	
WATER REACTIVITY	N	
RADIOACTIVITY	X	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	x	· · · · · · · · · · · · · · · · · · ·
рН	7N/A	
OXIDIZER	x	
PEROXIDE	x	
CYANIDE	x	
£ ∴FIDE	x	
CHLORIDE	x	
FL AMMA BLE	Y	Product burns
PCB	X	

2 0

Drum Id. : 8

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17H

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 50%

Drum Cond.: POOR

Overpack Size:

Jebris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	SLUDGE	BROWN	CLOUDY	100%
2	(Middle)				
3	(Bottom)				

'Manufacturer:

C_mical :

;enerator :

Misc. : Brown and white resin sludge with sand

2 1 9:23

Drum ID : 8

La er : (Top)

Physical State: SLUDGE

Color : BROWN

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	
WATER REACTIVITY	N	
RADIOACTIVITY	x	
WATER SOLUBILITY	Y	Partial
HEX/DCM SOLUBILITY	x	
рН	7	
OXIDIZER	x	
PEROXIDE	х	
CYANIDE	x	
. LFIDE	х	
CHLORIDE	х	
FLAMMABLE	х	
PCB	х	

2 2 3199

Drum Id. : 9

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time :

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17H

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 50%

Drum Cond.: POOR

Overpack Size: 85-GAL

Jebris/PPE:

No. of Layers: 2

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	riðnid	COLORLESS	CLEAR	50%
2	(Middle)				
3	(Bottom)	SLUDGE	GRAY	OPAQUE	50%

Manufacturer:

C mical :

Generator :

Misc. : Test positive for acetone Drager Tube

2 9 9000

Drum ID : 9

La er : (Top)

Color : COLORLESS

Physical State: LIQUID

Clarity : CLEAR

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	х	
WATER REACTIVITY	Y	color change to cloudy
RADIOACTIVITY	x	
WATER SOLUBILITY	Y	
HEX/DCM SOLUBILITY	x	
Н	7	
OXIDIZER	x	
PEROXIDE	x	
CYANIDE	x	
#FIDE	x	
CHLORIDE	Y	chloride risk
FLAMMABLE	N	
PCB	N	

Drum ID : 9

□ er : (Bottom)

Physical State: SLUDGE

Color : GRAY

Clarity : OPAQUE

2 2 3231

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	N	
WATER REACTIVITY	N	
RADIOACTIVITY	N	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	N	<i>;</i>
Нд	N	
OXIDIZER	N	
PEROXIDE	N	
CYANIDE	N	
LFIDE	N	
CHLORIDE	N	
FLAMMABLE	N	
PCB	N	

Drim Id. : 10

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 100%

Drum Cond.: FAIR

Overpack Size:

Debris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	SOLID	WHITE	CLOUDY	100%
2	(Middle)				
3	(Bottom)				

Manufacturer:

C. _mical

Generator :

Misc.

: writing on drum: 407# MCC 22-25

ADDITIONAL DRUM INFORMATION

Drum Class. :

Bulk Number :

Disposal Opt.:

Disposal Co. :

Manifest No. :

2 0 0203

rum ID : 10

T_K ≥r : (Top)

Color : WHITE

Physical State: SOLID

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	
WATER REACTIVITY	N	
RADIOACTIVITY	x	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	x	
рн	7	
OXIDIZER	х	
PEROXIDE	х	
CYANIDE	х	
LLFIDE	x	
CHLORIDE	x	
FLAMMABLE	N	
PCB	x	

2 2 3231

Drum Id. : 11

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 100%

Drum Cond.: FAIR

Overpack Size:

Debris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	SOLID	GRAY	CLOUDY	100%
2	(Middle)				
3	(Bottom)				

Manufacturer:

C. mical

Generator :

Misc.

2.0 0.05

rum ID : 11

Tai er : (Top)

Color : GRAY

Physical State: SOLID

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	
WATER REACTIVITY	N	
RADIOACTIVITY	х	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	х	
рн	7	
OXIDIZER	х	
PEROXIDE	x	
CYANIDE	х	
LFIDE	х	
CHLORIDE	x	
FLAMMABLE	Y	Sample "flashed" in ignitability test
PCB	х	

Drum Id. : 12

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 75%

Drum Cond.: POOR

Overpack Size:

Debris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	SOLID			
2	(Middle)				
3	(Bottom)				

Manufacturer:

C. nical :

Generator :

Misc. : Did not sample. Same as Drum #10.

rum ID : 12

T_{sk} ≥r : (Top)

Physical State: SOLID

Color :

Clarity :

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	N	
WATER REACTIVITY	N	
RADIOACTIVITY	N	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	N	
рН	N	
OXIDIZER	N	
PEROXIDE	N	
CYANIDE	N	
. LFIDE	N	
CHLORIDE	N	
FLAMMABLE	N	
PCB	N	

2 2 0203

Drum Id. : 13

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 75%

Drum Cond.: FAIR

Overpack Size: 85-GAL

Debris/PPE:

No. of Layers: 1

I	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)	SOLID	GREEN	CLOUDY	100%
2	(Middle)				
3	(Bottom)				

Manufacturer:

Chaical:

Generator :

Misc.

: Test positve for Styrene Drager Tube.

rum ID : 13

Tar : (Top)

Color : GREEN

Physical State: SOLID

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	
WATER REACTIVITY	N	
RADIOACTIVITY	х	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	х	
рH	7	
OXIDIZER	х	
PEROXIDE	x	
CYANIDE	х	
LFIDE	x	
CHLORIDE	х	
FLAMMABLE	x	
PCB	х	

2 1 02 0

Drum Id. : 14

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

Drum Top : BUNG

% Full : 100%

Drum Cond.: POOR

Overpack Size: 85-GAL

Debris/PPE:

No. of Layers: 1

LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1 (Top)	SOLID	GREEN	CLOUDY	100%
2 (Middle))	-		
3 (Bottom))			

Manufacturer:

C nical :

Generator

Misc. : Test positive for Styrene Drager Tube

2 2 9211

Drum ID : 14

L er : (Top)

Physical State: SOLID

Color : GREEN

TEST NAME	RESULT	COMMENTS
AIR REACTIVITY	x	
WATER REACTIVITY	N	
	x	
RADIOACTIVITY	ĺ	
WATER SOLUBILITY	N	
HEX/DCM SOLUBILITY	X	
рН	7	
OXIDIZER	x	
PEROXIDE	x	
CYANIDE	x	
LFIDE	x	
CHLORIDE	х	
FLAMMABLE	N	
PCB	x	

2 1 1212

Drum Id. : 15

Location: pit

Date

: 01/10/95

Sampler : S.Getty\M.Morganti

Time :

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size :

Drum Top : BUNG

% Full

Drum Cond.: FAIR

Overpack Size:

Debris/PPE:

No. of Layers:

LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1 (Top)				
2 (Middle)				
3 (Bottom)				

Manufacturer:

C. nical :

Generator

Misc.

: Rock solid to the top. Could not even open. Not sampled.

Drum Id. : 16

Location: pit

)are

: 01/10/95

Sampler : S.Getty\M.Morganti

Time

Witness : S.Getty\M.Morganti

Drum Type : STEEL-17E

Drum Size : 55-GAL

rum Top : BUNG

% Full : 100%

Drum Cond.: FAIR

Overpack Size: 85-GAL

Jebris/PPE:

No. of Layers: 1

	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)				
2	(Middle)				
3	(Bottom)				

'Manufacturer:

c mical :

Generator :

Misc. : Rock solid to the top. Not sampled.

2 2 3214

Drum Id. : 1,8,9,11 C

Location: composite

Date

: 01/19/95

Sampler : S.Getty

Time

.

Witness:

Drum Type :

Drum Size :

Drum Top :

% Full

Drum Cond.:

Overpack Size:

Debris/PPE:

No. of Layers:

LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1 (Top)				
2 (Middle)				
3 (Bottom)				

Manufacturer:

C. nical :

Generator

Misc.

: Sample is a composite of 4 Drum samples (#1,8,9, and 11).

Drum Id. : COMPOSITE2

Location:

)ate

: 01/19/95

Sampler : S.Getty

Time

Witness :

Drum Type :

Drum Size :

rum Top :

% Full :

Drum Cond.:

Overpack Size:

Jebris/PPE:

No. of Layers:

== 	LAYERS	PHYSICAL STATE	COLOR	CLARITY	LAYER DEPTH
1	(Top)				
2	(Middle)				
3	(Bottom)				

Manufacturer:

C mical :

Senerator :

Misc. : Sample is a composite of samples 3,7,10,13, and 14.

Appendix C

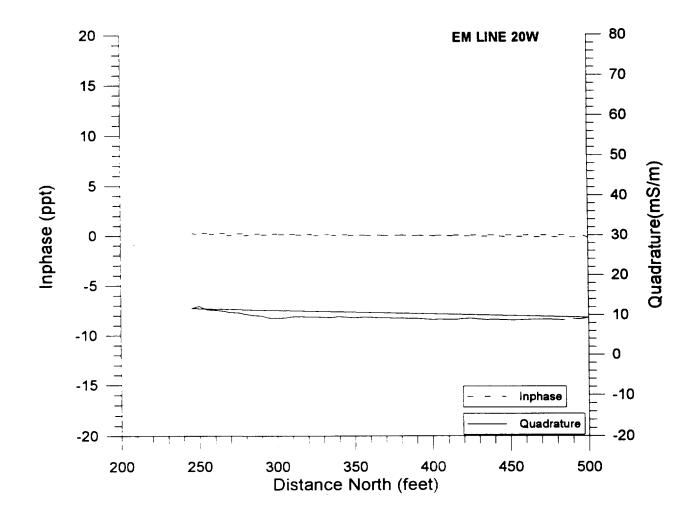
2 1 3217

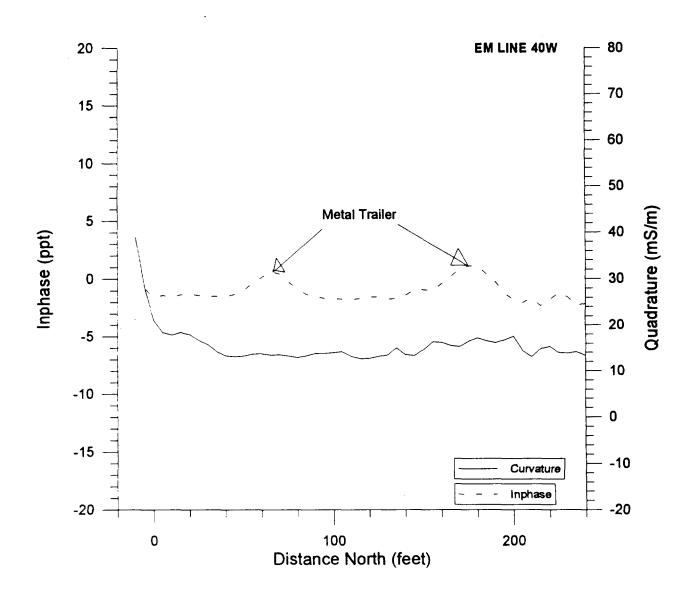
APPENDIX C

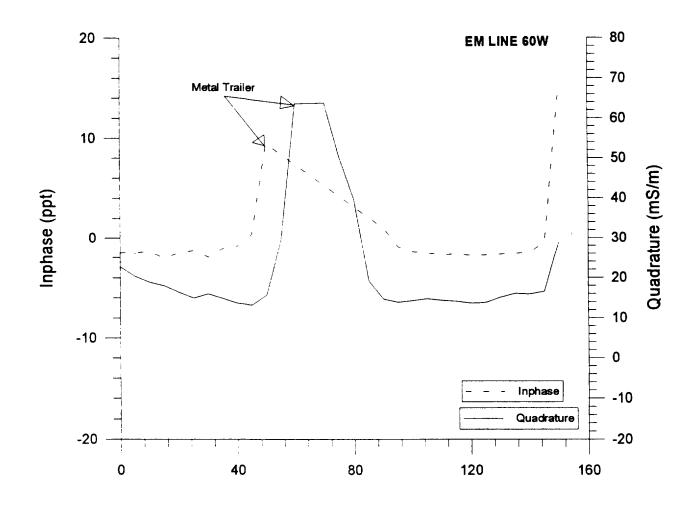
GEOPHYSICAL SURVEY
Pier Drum Site
Bradenton, FL
January 1995

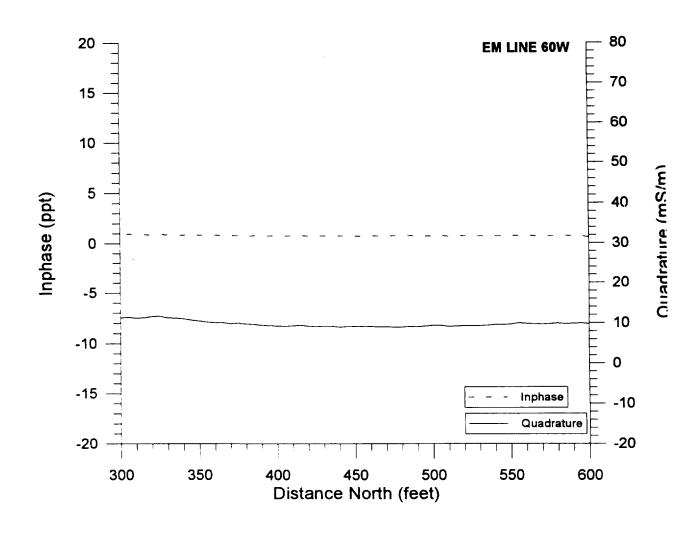
Appendix Contents:

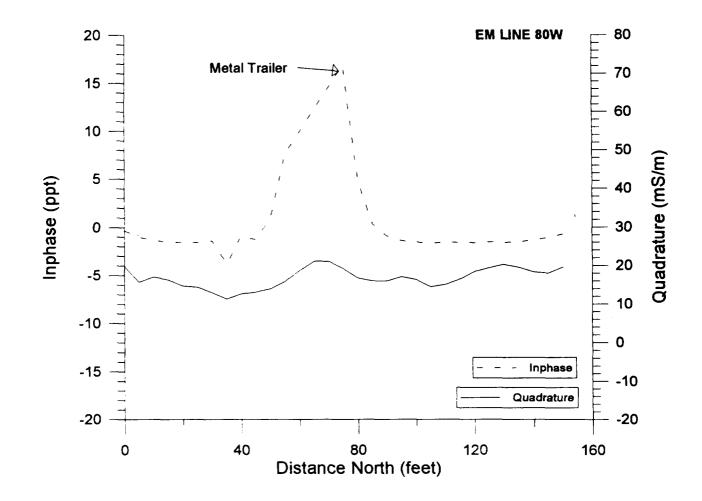
Electromagnetic (EM-31) Profiles Magnetic Profiles

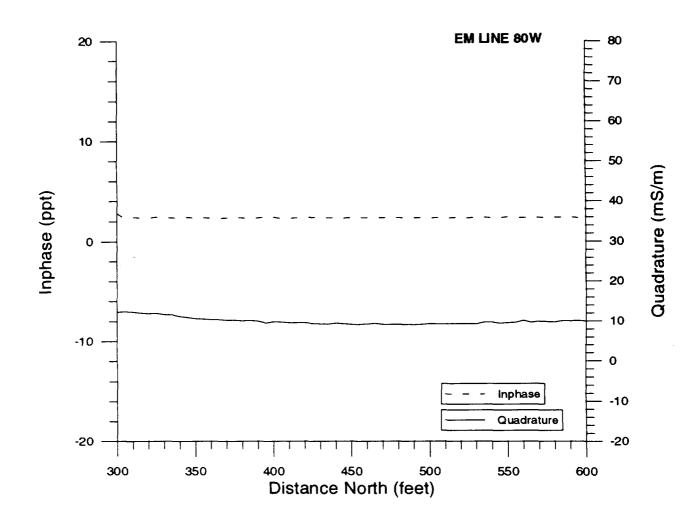


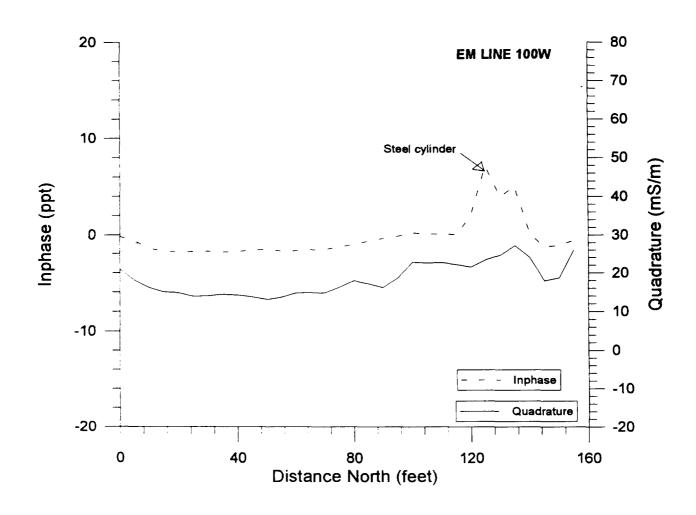


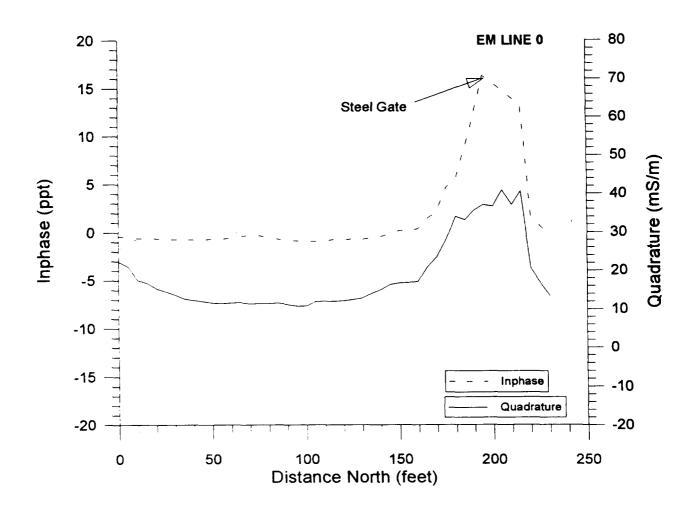


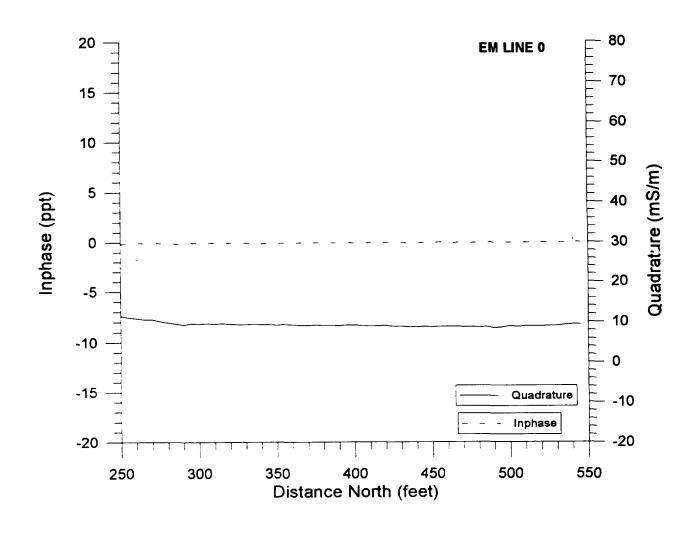


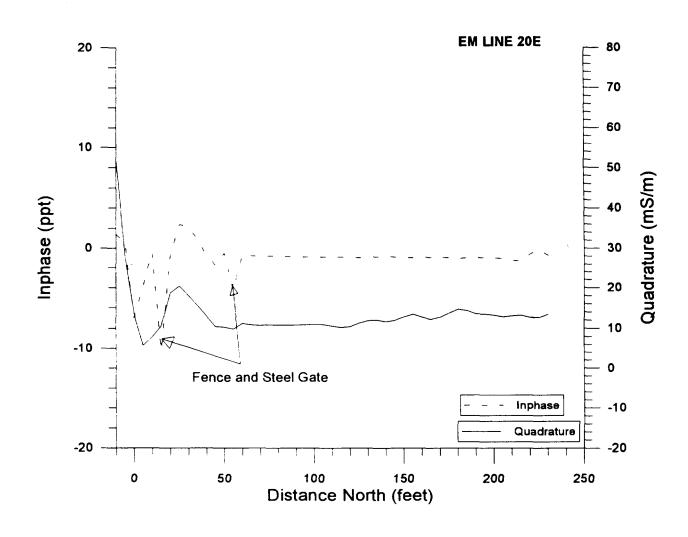


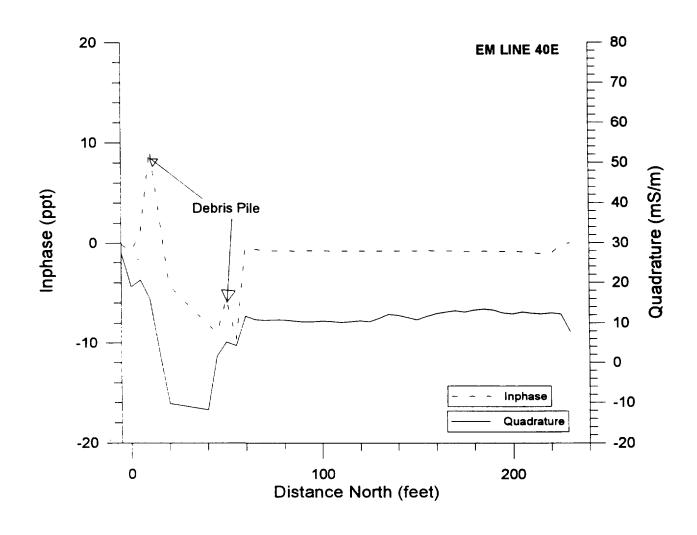


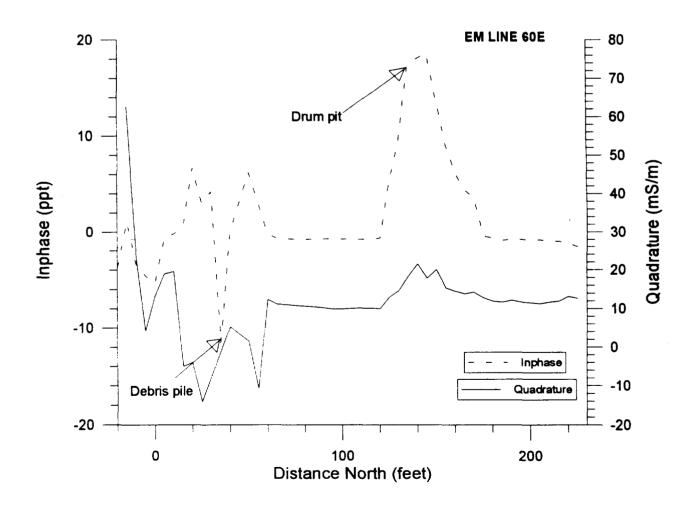


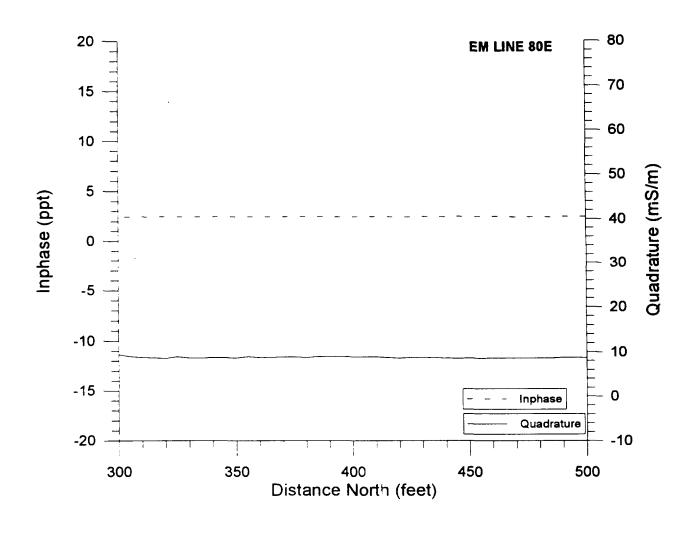


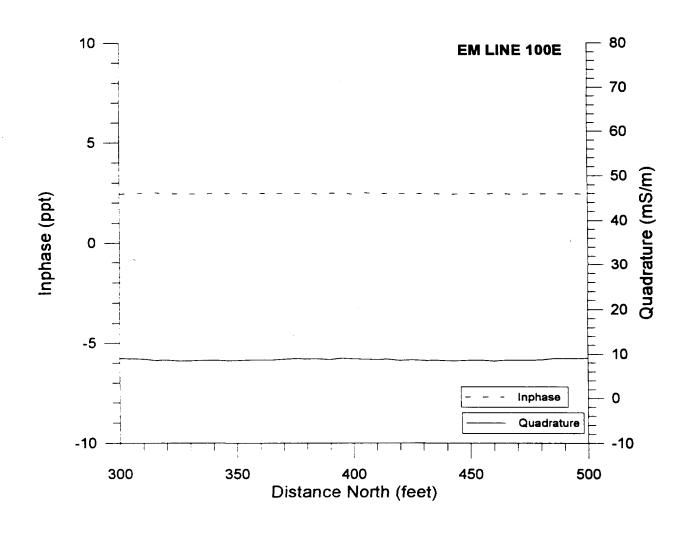


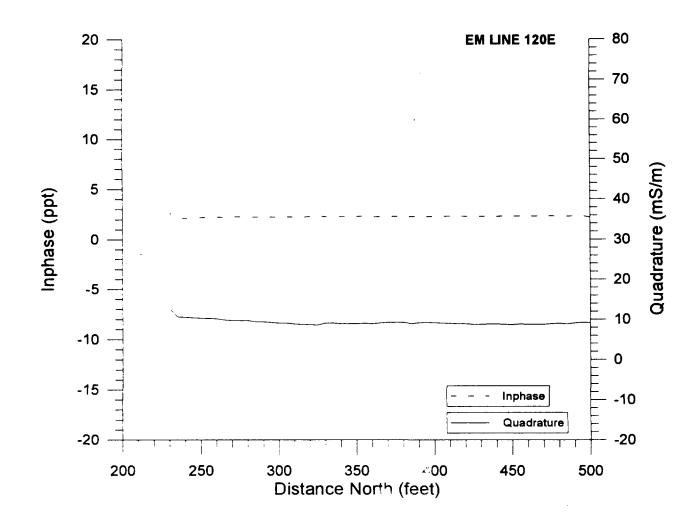


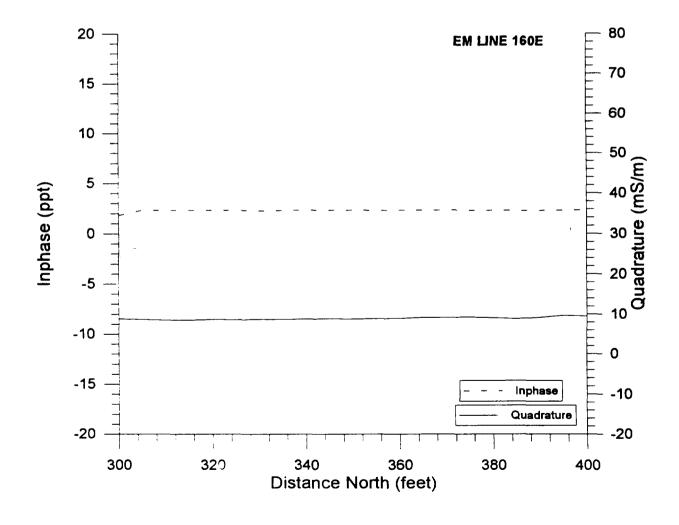


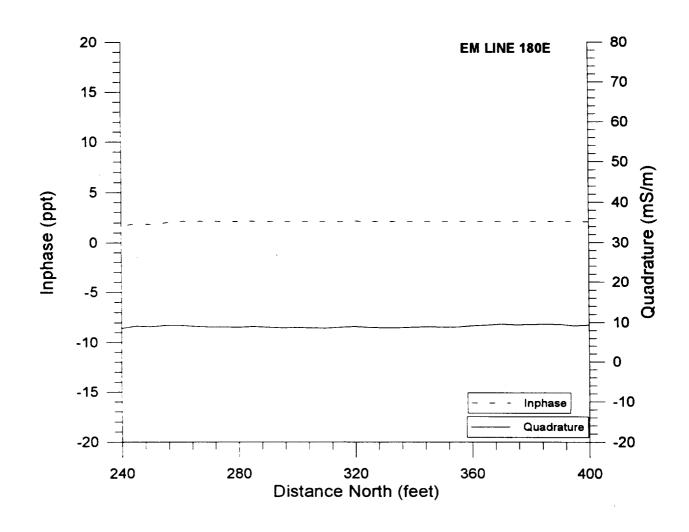


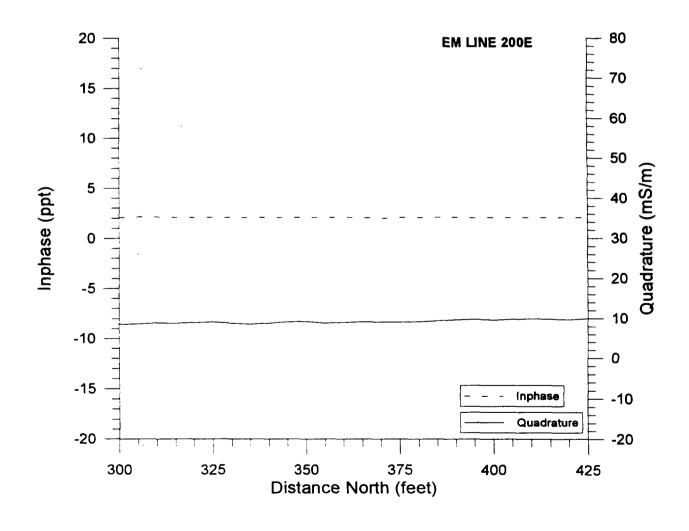


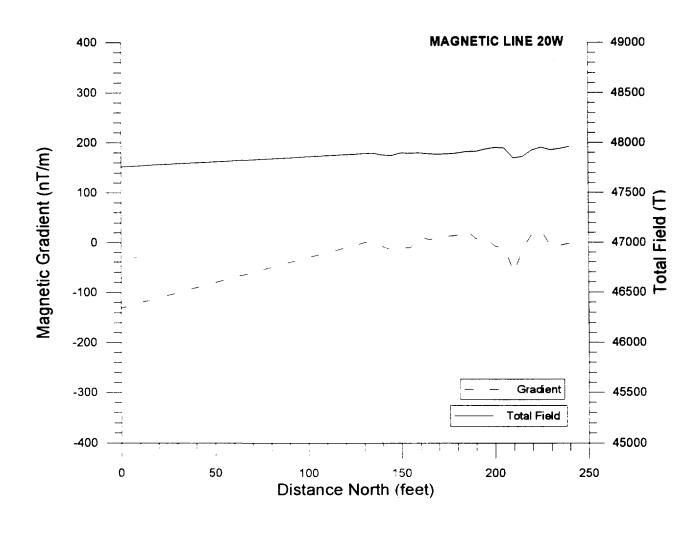


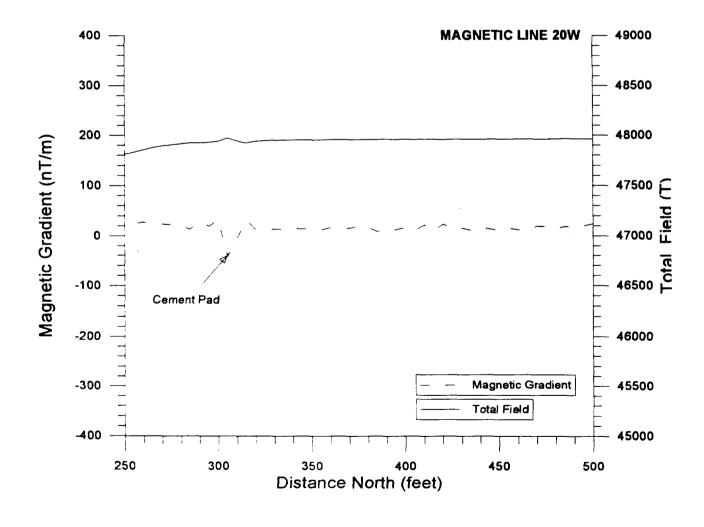


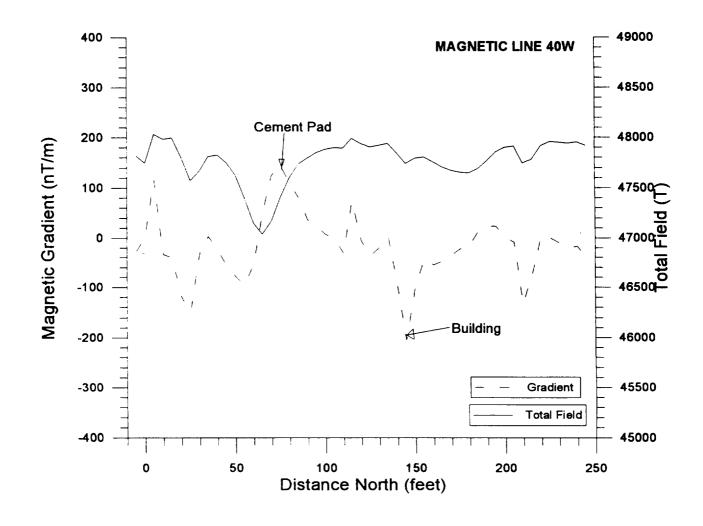


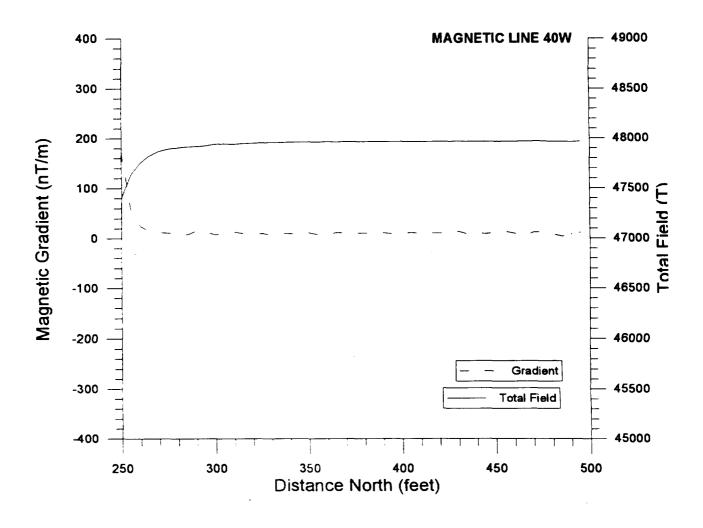


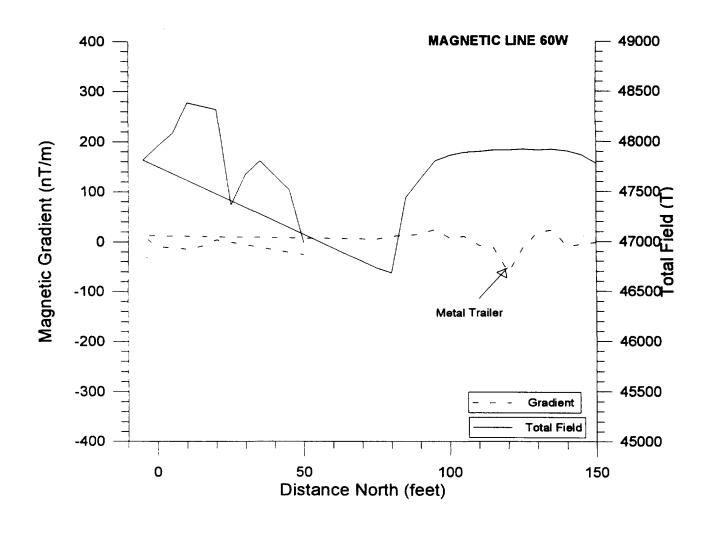


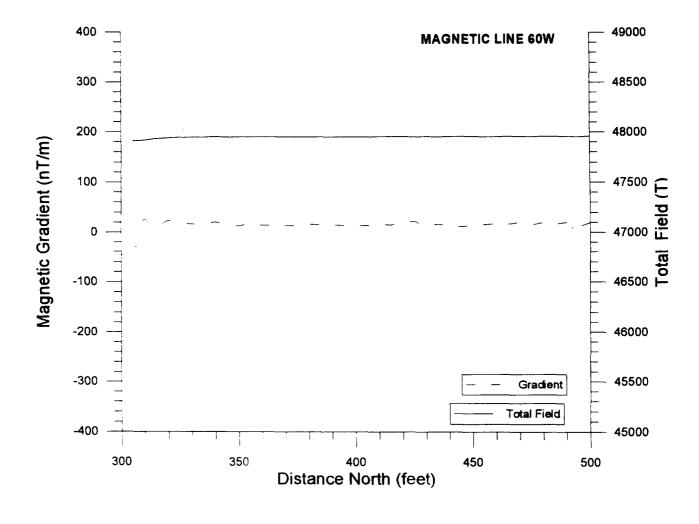


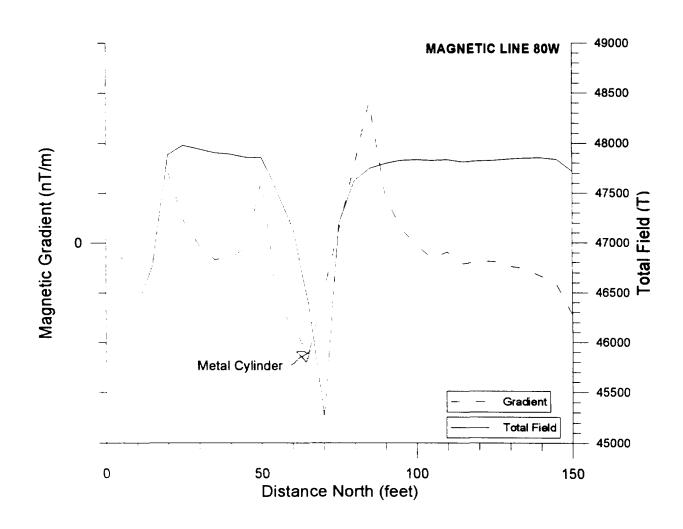


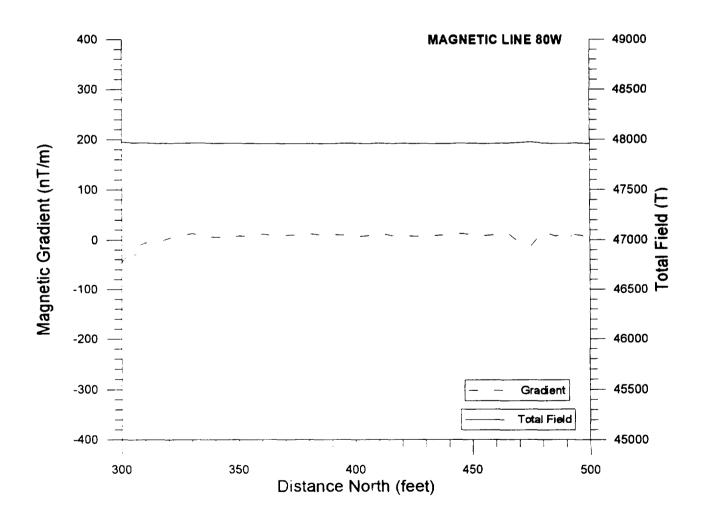


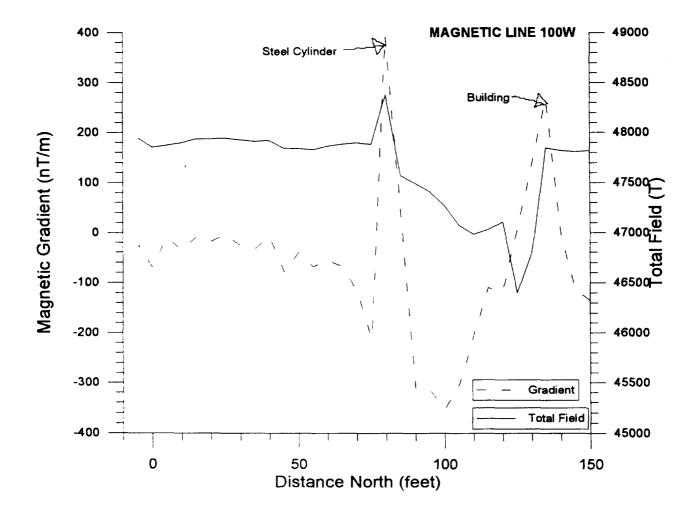


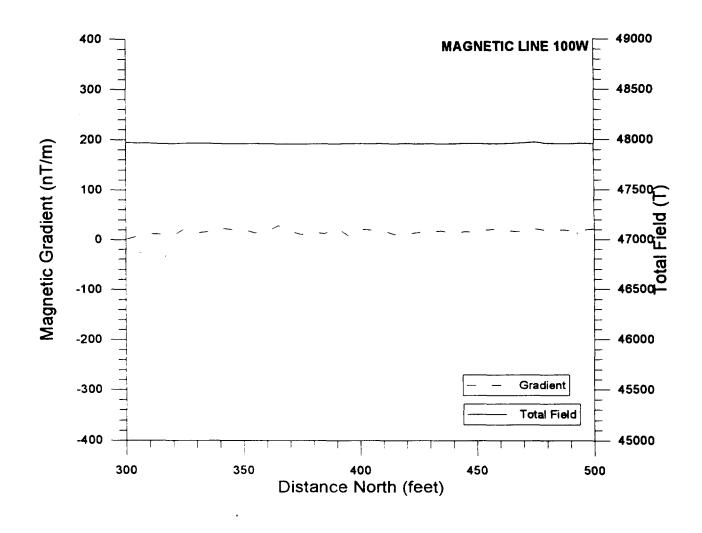


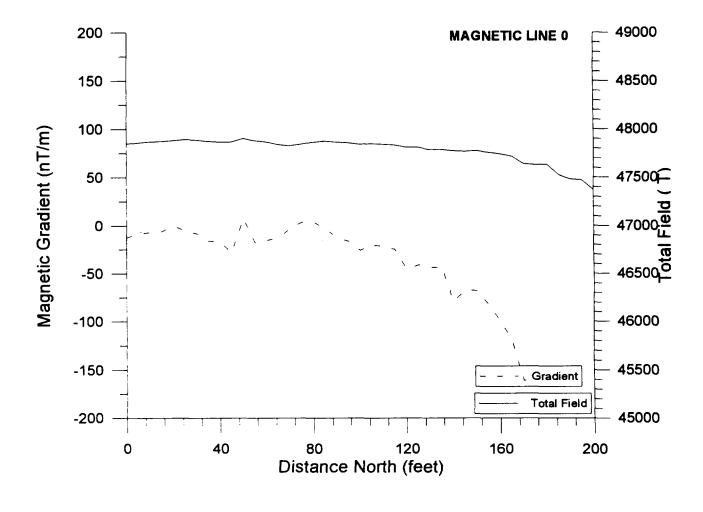


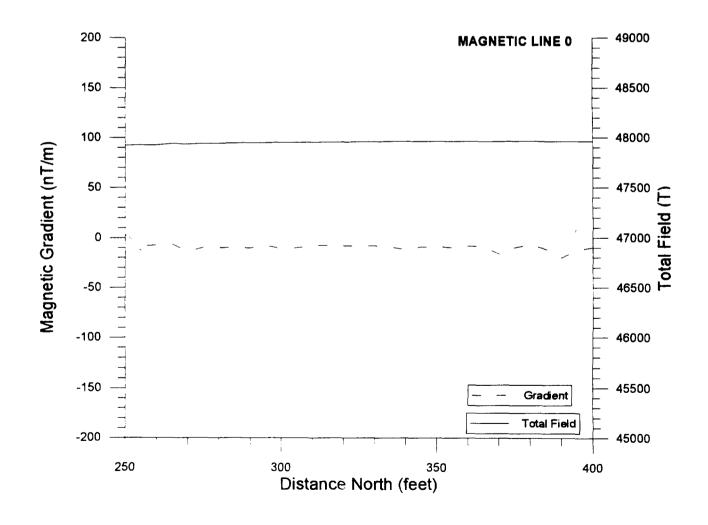


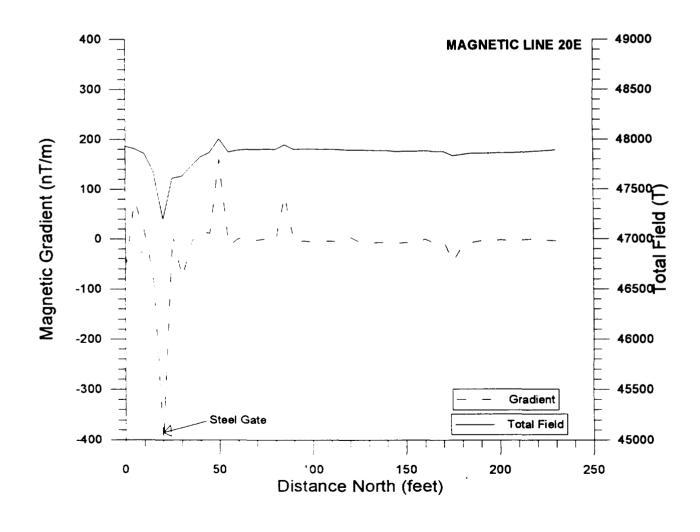


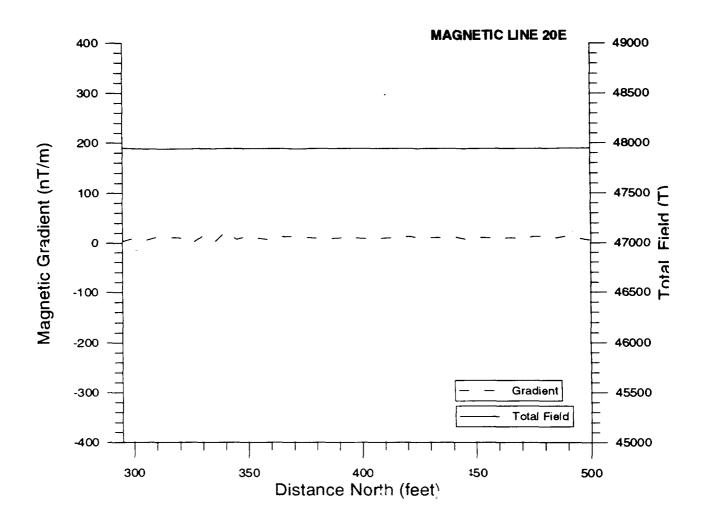


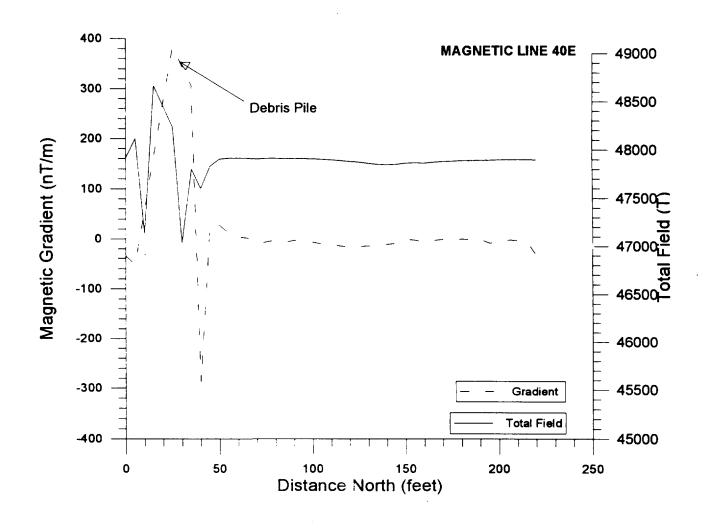


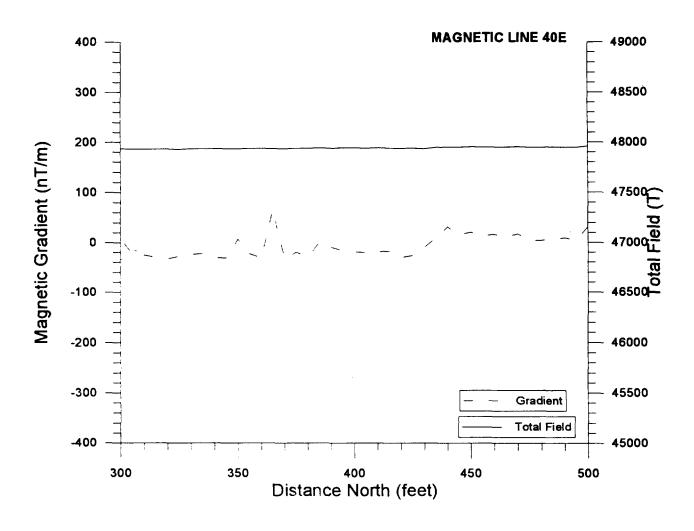


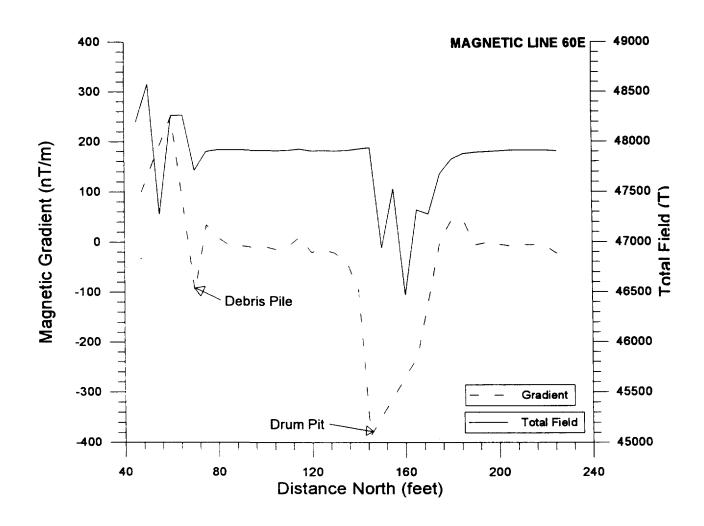


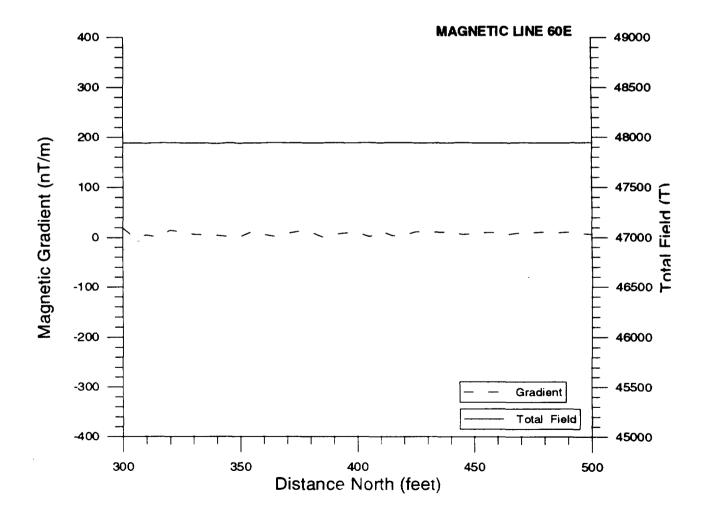


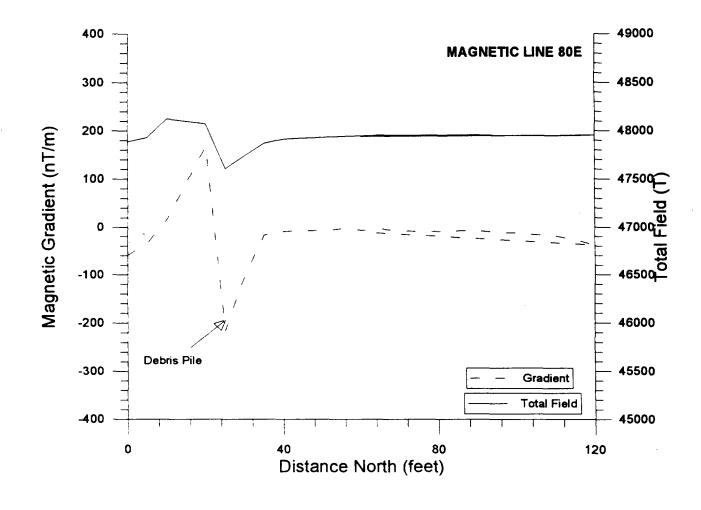


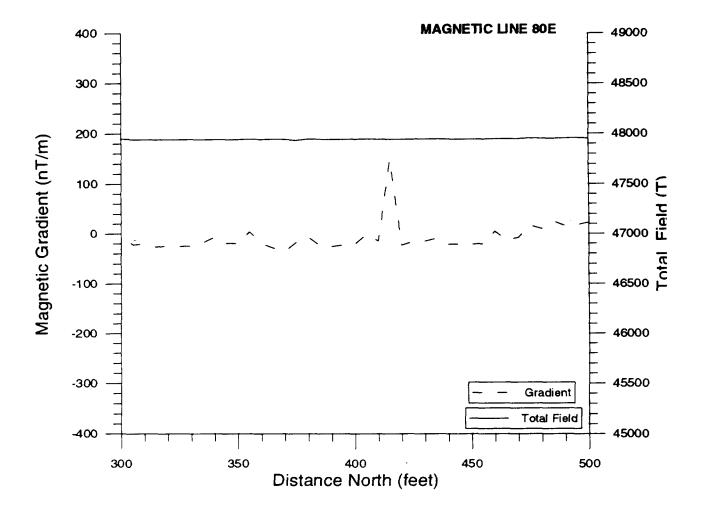






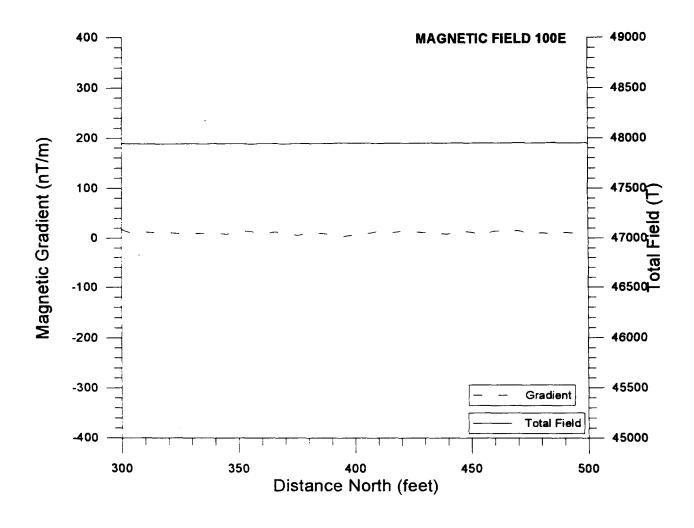


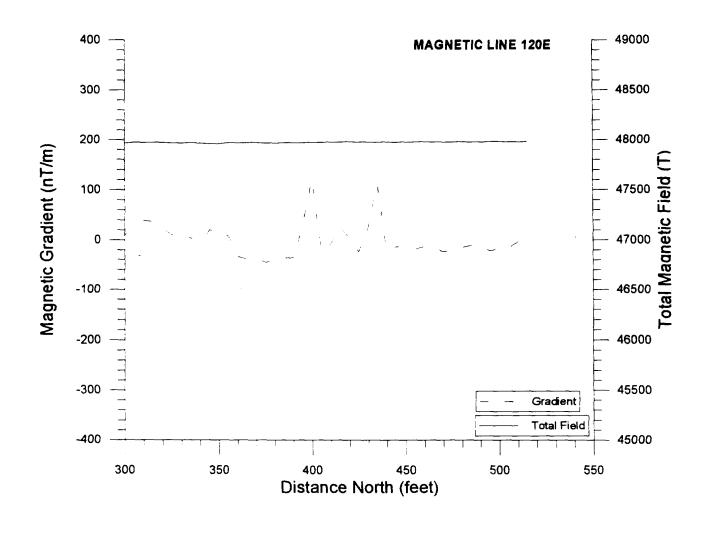


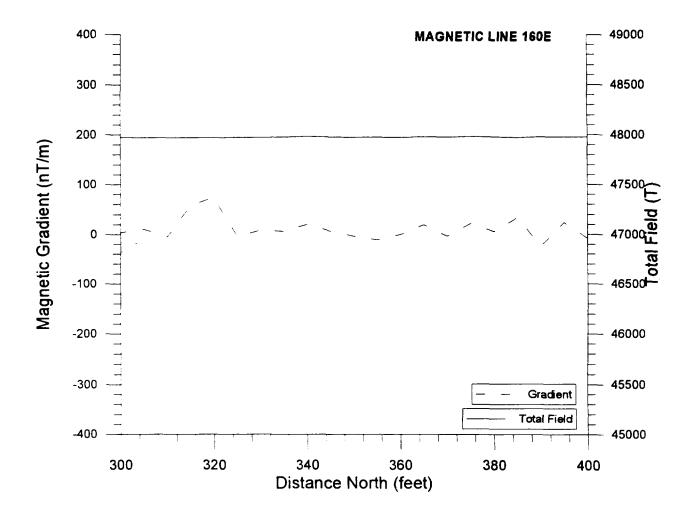


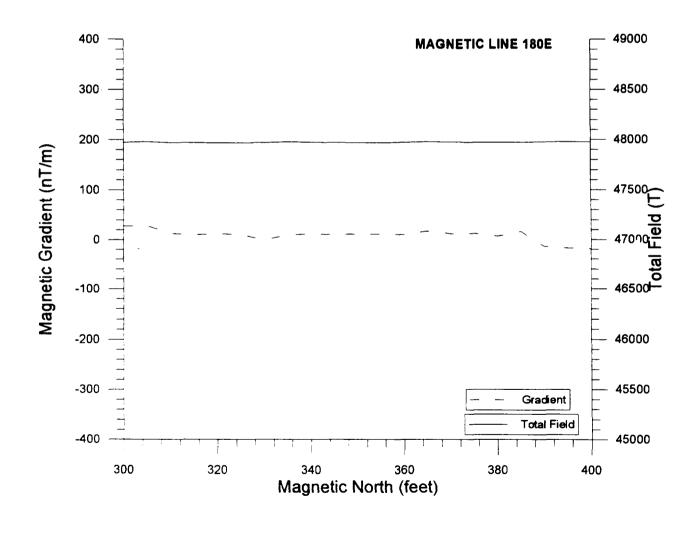
,

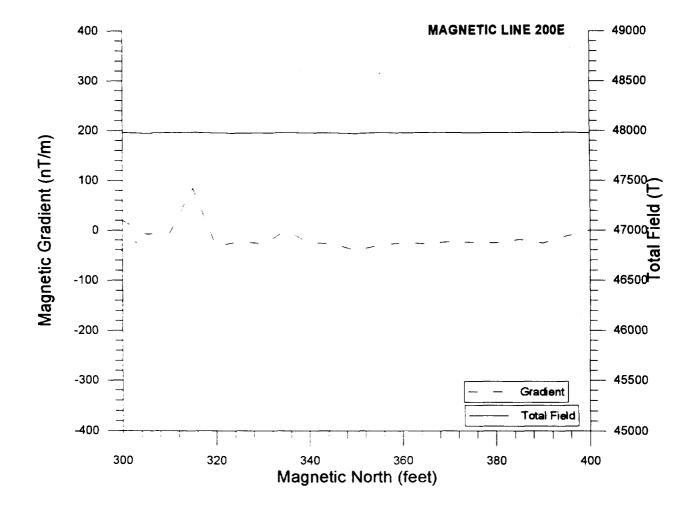
* 3











Appendix D

2 2 22 22 2

APPENDIX D

Final Analytical Report Pier Drum Site Bradenton, FL January 1995

Appendix Contents:

Analytical Procedures and Results Quality Assurance/Quality Control Chain of Custody Communications





Roy F. Weston, Inc. GSA Raritan Depot Building 209 Annex (Bay F) 2890 Woodbridge Avenue Edison, New Jersey 08837-3679 908-321-4200 • Fax 908-494-4021

DATE:

March 21, 1995

TO:

R. Singhvi

EPA/ERT

FROM:

V. Kansal

Analytical Section Leader Vinor Kausal

SUBJECT: DOCUMENT TRANSMITTAL UNDER WORK ASSIGNMENT # 0-098

Attached please find the following document prepared under this work assignment:

Pier Drum - Analytical Report

Central File WA # 0-098

(w/attachment)

R. Lewis

Task Leader

G. Armstrong

Data Validation and Report Writing Group Leader

G. Powell

Work Assignment Manager

ANALYTICAL REPORT

Prepared by Roy F. Weston, Inc.

> Pier Drum Bradenton, FL

March 1995

EPA Work Assignment No. 0-098 WESTON Work Order No. 03347-040-001-0098-01 EPA Contract No. 68-C4-0022

> Submitted to G. Powell **EPA-ERT**

Task Leader

Analysis by:

REAC

Analab

Prepared by:

G. Karustis

Ralph M. Shapot

Analytical Section Leader

Program Manager

Reviewed by:

G. Armstrong

Table of Contents

Topic Page Numbers Introduction Page 1 Case Narrative Page 2 Summary of Abbreviations Page 9 Section I **Analytical Procedures** Analytical Procedure for VOA in Water and Drum Liquids Page 10 Analytical Procedure for VOA in Soil Page 12 Analytical Procedure for TCLP VOA Page 14 Analytical Procedure for BNA in Water, TCLP Extracts and Drum Liquids Page 15 Analytical Procedure for BNA in Soil Page 17 Analytical Procedure for TCLP BNA Page 19 Analytical Procedure for Pesticides/PCBs in Water 20 Page Analytical Procedure for Pesticides/PCBs in Soil 22 Page Analytical Procedure for Pesticides/PCBs in Drum Liquids Page 24 Analytical Procedure for TCLP Pesticides/PCBs Page 26 Analytical Procedure for Metals in Water Page 26 Analytical Procedure for Metals in Soil Page 26 Analytical Procedure for TCLP Metals Page 27 Analytical Procedure for Metals in Drum Liquid Samples Page 27 Analytical Procedure for TCLP Metals in Drum Solids Page 27 Analytical Procedure for BTU Page 27 Analytical Procedure for Ash Page 27 Analytical Procedure for Chlorine Page 28 Analytical Procedure for Sulfur Page 28 Analytical Procedure for Moisture 28 Page **VOA Results** Results of the VOA Analysis for the Water Samples Table 1.1 Page 29 Results of the VOA in Water TICs Table 1.2 Page 39 Results of the VOA Analysis for the Soil Samples Table 1.3 Page 40 Results of the VOA in Soil TICs Table 1.4 Page 44 Results of the TCLP VOA Analysis Table 1.5 45 Page Results of the VOA Analysis for the Drum Liquid Samples Table 1.6 Page 46 Results of the VOA in Drum Liquid TICs Table 1.7 50 Page Results of the TCLP WOA Analysis for the Drum Solid Samples Table 1.8 Page 51 **BNA Results** Results of the BNA in Water Analysis Table 1.9 Page 52 Results of Tentatively Identified Compounds for BNA in Water **Table 1.10** Page 56 Results of the BNA in Soil Analysis Table 1.11 Page 65 Results of Tentatively Identified Compounds for BNA in Soil Table 1.12 Page 67 Results of the TCLP BNA Analysis Table 1.13 71 Page Results of Tentatively Identified Compounds for BNA in TCLP Table 1.14 Page 73 Results of the BNA in Drum Liquid Analysis Table 1.15 77 Page Results of Tentatively Identified Compounds for BNA in Drum Liquid Table 1.16 Page 81 Results of the BNA Analysis for the TCLP Extracts from the Drum Solids

Table 1.17

86

Page

	Pesticide\PCB Results			
	Results of the Pesticides/PCB Analysis for the Water Samples	Table 1.18	Page	87
	Results of the Pesticides/PCB Analysis for the Soil Samples	Table 1.19	Page	89
	Results of the TCLP Pest/PCB Analysis	Table 1.20	Page	90
	Results of the Pest/PCB Analysis in Drum Liquid Samples	Table 1.21	Page	91
	Results of the Pest/PCB Analysis in Drum Solid Samples	Table 1.22	Page	92
	Metals Results			
	Results of the Metals Analysis for the Water Samples	Table 1.23	Page	93
	Results of the Metals Analysis for the Soil Samples	Table 1.24	Page	96
	Results of the TCLP Metals Analysis	Table 1.25	Page	97
	Results of the Metals Analysis for the Drum Liquid Samples	Table 1.26	Page	98
	Results of the Metals Analysis in TCLP			
	Extracts from the Drum Solids	Table 1.27	Page	99
	Wet Chemistry Results			
	Results of the Wet Chemistry Analysis	Table 1.28	Page	100
Section	II			
	QA/QC for VOA		Daga	101
	Results of the Internal Standard Areas and Surrogate Recoveri	AC	Page	101
	for the Water Samples	Table 2.1	Page	103
	Results of the MS/MSD Analysis for the Water Samples	Table 2.1	Page	103
	Results of the Internal Standard Areas and Surrogate Recoverience		rage	104
	for the Soil Samples	Table 2.3	Page	105
	Results of the MS/MSD Analysis for the Soil Samples	Table 2.4	Page	106
	Results of the Internal Standard Areas and Surrogate Recoverience		1 age	100
	for the Drum Liquid Samples	Table 2.5	Page	107
	Results of the MS/MSD Analysis for the Drum Liquid Samples		Page	108
	Results of the Initial Calibration	Table 2.7	Page	109
	Results of the Continuing Calibration	Table 2.8	Page	115
	Correlation of Samples and Calibrations	Table 2.9	Page	127
	Results of the TCLP Surrogate Recoveries	Table 2.10	Page	128
	Results of the TCLP Internal Standard Areas	Table 2.11	Page	129
	Results of the MS/MSD Analysis for the TCLP VOA Samples	Table 2.12	Page	130
	Results of the TCLP Surrogate Recoveries	I doic 2.12	rage	150
	for the Drum Solid Samples	Table 2.13	Page	131
	Results of the TCLP Internal Standard Areas	14010 2.13	r agc	131
	for the Drum Solid Samples	Table 2.14	Page	132
	Results of the Blank Spike Analysis for the TCLP Analysis	14010 2.14	1 age	132
	of Drum Solid Samples	Table 2.15	Page	133
	Results of the MS/MSD Analysis for the TCLP VOA Analysis	ravic 2.13	1 age	133
	of the Drum Solid Samples	Table 2.16	Page	134
	of the Diam Solid Samples	1 aut 2.10	ı agc	134

QA/QC for BNA			Page	135
Results of the Internal Standard Areas and Surrogate Recoveries				
for the Water Samples	Table	2.17	Page	137
Results of the MS/MSD Analysis for the Water Samples	Table	2.18	Page	138
Results of the Internal Standard Areas and Surrogate Recoveries				
for the Soil Samples	Table	2.19	Page	139
Results of the MS/MSD Analysis for the Soil Samples	Table	2.20	Page	140
Results of the Internal Standard Areas and Surrogate Recoveries			_	
for the TCLP Samples	Table	2.21	Page	141
Results of the MS/MSD Analysis for the TCLP Samples	Table	2.22	Page	142
Results of the Internal Standard Areas and Surrogate Recoveries				
for the Drum Liquid Samples	Table	2.23	Page	143
Results of the MS/MSD Analysis for the Drum Liquid Samples	Table	2.24	Page	144
Results of the Initial Calibration	Table	2.25	Page	145
Results of the Continuing Calibration	Table	2.26	Page	147
Correlation of Samples and Calibrations	Table	2.27	Page	161
Results of the TCLP Surrogate Recoveries for the Drum Solid Sa	mples			
	Table	2.28	Page	162
Results of the TCLP Internal Standard Areas for the Drum Solid	Sampl	es		
	Table	2.29	Page	163
Results of the Blank Spike Analysis for the TCLP BNA Analysis	Table	2.30	Page	164
Results of the MS/MSD Analysis for the TCLP BNA Analysis				
of the Drum Solid Samples	Table	2.31	Page	165
QA/QC for Pesticides/PCBs			Page	166
Results of the Surrogate Recoveries in Water	Table	2.32	Page	168
Results of the MS/MSD Analysis for Pest/PCB in Water	Table		Page	169
Results of the Surrogate Recoveries in Soil	Table	2.34	Page	170
Results of the MS/MSD Analysis for Pest/PCB in Soil	Table		Page	171
Results of the Surrogate Recoveries in TCLP	Table		Page	172
Results of the MS/MSD Analysis for Pest/PCB in TCLP	Table		Page	173
Results of the Surrogate Recoveries in Drum Liquids	Table		Page	174
Results of the MS/MSD Analysis for Pest/PCB in Drum Liquids	Table		Page	175
Results of the Surrogate Recoveries for the Drum Solid Samples		2.40A	Page	176
Results of the Surrogate Recoveries for the Drum Solid Samples		2.40B	Page	176
Results of the Blank Spike Analysis for the Drum Solid Samples	Table		Page	177
Results of the MS/MSD Analysis for the Drum Solid Samples	Table	2.42	Page	178
QA/QC for Metals		2 42	Page	179
Results of the QC Standard Analysis for Water Samples	Table		Page	181
Results of the MS/MSD Analysis for Water Samples	Table		Page	182
Results of the Blank Spike Analysis for Water Samples	Table		Page	183
Results of the QC Standard Analysis for Soil Samples	Table		Page	184
Results of the MS/MSD Analysis for Soil Samples	Table		Page	185
Results of the Blank Spike Analysis for Soil Samples	Table		Page	186
Results of the QC Standard Analysis for TCLP Samples	Table		Page	187
Results of the MS/MSD Analysis for TCLP Samples	Table		Page	188
Results of the OC Standard Applysis for TCLP Samples	Table		Page	189
Results of the QC Standard Analysis for Drum Liquid Samples	Table		Page	190
Results of the Blank Spike Analysis for Drum Liquid Samples	Table		Page	191
Results of the Blank Spike Analysis for Drum Solid Samples Results of the MS/MSD yealysis for the TCLP Analysis	Table	4.34	Page	192
Results of the MS/MSD analysis for the TCLP Analysis of the Drum Solid Samples	Table	2 55	Daga	102
of the Diam Sond Samples	I auic	در. ے	Page	193

QA/QC for BTU, Ash, Moisture, Sulfur and Chlorine 195 Page Results of the Duplicate Analysis for the Drum Liquid Samples Table 2.56 Page 196 Section III Chains of Custody Page 197 Communications Page 204 **Appendices** Appendix A Data for VOA In Water and Sediment Page E 23 01 For TCLP Page E 47 01 In Drum Liquids Page E 32 01 Appendix B Data for BNA In Water and Sediment Page E 33 01 In Drum Liquids Page E 67 01 Appendix C Data for Pest/PCB In Water Page E 42 01 In Soil Page E 51 01 In Drum Liquids Page E 58 01 Appendix D Data for Metals In Water Page E 30 01 In Soil Page E 35 01 Page E 52 01 For TCLP In Drum Liquids Page E 56 01 Appendix E Data for Drum Solids and Wet Chemistry Page E 78 01

Appendices will be furnished on request.

INTRODUCTION

REAC Laboratory, in response to WA # 0-098, provided analytical support for water, soil, drum solid and drum liquid samples collected from the Pier Drum Site in Bradenton, FL. This support included the analysis of water, soil and drum liquid samples as described in the following table. The support also included the subcontracted analysis of drum solid and drum liquid samples, QA/QC, data review, and preparation of an analytical report containing a summary of the analytical methods, the results, and the QA/QC results.

The samples were treated with procedures consistent with those described in SOP # 1008 and are summarized in the following table:

COC #**	Number of Samples	Sampling Date	Date Received	Matrix	Analysis	Laboratory
00065	3	1/10/95	1/30/95	Drum Liquid	VOA, BNA, Pest/PCB, Metals	REAC
00066	3	1/10/95	1/30/95	Drum Solid	TCLP VOA, TCLP BNA, TCLP Pest/PCB, TCLP Metals	Analab
00066	3	1/10/95	1/30/95	Drum Liquid	%Cl, Moisture, BTU, Ash, %S	Wastex
03663	1	1/11/95	1/13/95	Water	BNA Pest/PCB, TAL Metals	REAC
03663	2	1/11/95	1/13/95	Water	VOA	REAC
03664	3	1/11/95	1/13/95	Soil	TCLP VOA	Analab
03670	3	1/11/95	1/13/95	Water	Water VOA, BNA, Pest/PCB, TAL Metals	

^{**} COC # denotes Chain of Custody number

The sample table is continued on the next page

2 1 0210

Sample Table (Cont)

COC #**	Number of Samples	Sampling Date	Date Received	Matrix	Analysis	Laboratory
03671	2	1/11/95	1/13/95	Water	VOA, BNA, Pest/PCB, TAL Metals	REAC
03672	2	1/11/95	1/13/95	Water	VOA, BNA, Pest/PCB, TAL Metals	REAC
03672	3	1/11/95	1/13/95	Soil	VOA, BNA, Pest/PCB, TAL Metals	REAC
10027	3	1/18/95	1/18/95	Soil	TCLP Metals, TCLP BNA, TCLP Pest/PCB	REAC

^{**} COC # denotes Chain of Custody number

CASE NARRATIVE

VOA Package E 32

These samples are drum liquids and the holding time criterion does not apply. The analysis was conducted to determine the means of disposal of the drum contents.

VOA Package E 23

The continuing calibration run on 1/13/95 exceeds the 25% difference requirement for acetone (46.24%) and 2-hexanone (27.32%). These analytes were not detected in the samples run under this calibration; the data are not affected.

The continuing calibration run on 1/14/95 exceeds the 25% difference requirement for acetone (29.70%). This analyte was not detected in the samples run under this calibration; the data are not affected.

The continuing calibration run on 1/17/95 exceeds the 25% difference requirement for acetone (52.75%). Both detected and non detected values will be considered as estimated for Lab blank - 1/17/95 and sample 00642.

TCLP VOA Package E 47

The method blank analyzed on 1/20/95 contained 2.6 ug/l methylene chloride. None of the associated samples contained methylene chloride. The data are not affected.

The continuing calibration run on 1/20/95 exceeds the 25% difference requirement for chloromethane (35.08%), bromomethane (27.44%), chloroethane (30.66%), acetone (36.90%), 2-butanone (31.26%), bromoform (32.68%), 4-methyl-2-pentanone (34.49%) and 2-hexanone (35.93%). These analytes were not detected in the samples quantitated by this continuing calibration. The data are not affected.

One surrogate percent recovery (1,2-dichloroethane-d₄) exceeds the acceptable QC limits for sample B 00644. All results for this sample should be considered as estimated.

BNA Package E 33

The method blank analyzed on 1/27/95 contained 46 ug/l di-n-butyl phthalate, 5 ug/l butyl benzyl phthalate and 11 ug/l bis(2-ethylhexyl)phthalate. The positive results of less than 230 ug/l di-n-butyl phthalate, 25 ug/l butyl benzyl phthalate and 55 ug/l bis(2-ethylhexyl)phthalate in samples B 00525, B 00526 and B 00527 are considered to be non-detected.

The method blank analyzed on 1/30/95 contained 2 ug/l di-n-butyl phthalate and 1 ug/l bis(2-ethylhexyl)phthalate. The positive results of less than 10 ug/l di-n-butyl phthalate and 5 ug/l bis(2-ethylhexyl)phthalate in samples CD 00636, CD 00638, CD 00635, CD 00640 and CD 00634 are considered to be non-detected.

The method blank analyzed on 1/31/95 contained 2400 ug/kg di-n-butyl phthalate, 150 ug/kg butyl benzyl phthalate and 620 ug/kg bis(2-ethylhexyl)phthalate. The positive results of less than 12000 ug/kg di-n-butyl phthalate, 750 ug/kg butyl benzyl phthalate and 3100 ug/kg bis(2-ethylhexyl)phthalate in samples A 00642, A 00643 and A 00644 are considered to be non-detected.

The initial calibration run on 11/4/94 exceeds the acceptable QC limits for diethyl phthalate (31.079%). Since no samples were quantitated by this initial calibration, the data are not affected.

The continuing calibration run on 1/27/95 exceeds the acceptable QC limits for hexachlorobenzene (26.92%). Since this analyte was not detected in any of the samples, the data are not affected.

The continuing calibration run on 1/30/95 exceeds the acceptable QC limits for hexachlorobutadiene (26.99%) and hexachloropentadiene (34.84%). Since these analytes were not detected in any of the samples, the data are not affected.

The continuing calibration run on 1/31/95 exceeds the acceptable QC limits for hexachlorobutadiene (34.48%). Since this analyte was not detected in any of the samples, the data are not affected.

2 2 32/2

The continuing calibration run on 2/1/95 exceeds the acceptable QC limits for hexachlorobutadiene (38.75%). Since this analyte was not detected in any of the samples, the data are not affected.

Two base/neutral surrogate recoveries exceed the acceptable QC limits for nitrobenzene-d₅ (27%) and 2-fluorobiphenyl (30%) for WBLK 011895. The positive base/neutral results detected in this blank should be regarded as estimated.

One acid surrogate recovery exceeds the acceptable QC limits for 2,4,6-tribromophenol (125%) for sample CD 00638. The data are not affected.

For sample BE 00527 one relative percent difference for 1,4-dichlorobenzene (29%) and three spike percent recoveries (MS and MSD for 4-nitrophenol and MSD for pentachlorophenol) exceeded the acceptable QC limits. The data are not affected.

For sample CD 00644 one spike percent recovery (4-nitrophenol) exceeded the acceptable QC limits. The data are not affected.

For sample A 00644 three spike percent recoveries (MS for phenol and 2,4-dinitrotoluene and the MSD for 2,4-dinitrophenol) exceeded the acceptable QC limits. The data are not affected.

BNA Package E 67 (Drum Liquids

These samples are drum liquids and the holding time criterion is not applicable.

The bottom layer of sample 00913 was diluted and an aliquot was analyzed. Surrogates were not added and an MS/MSD analysis was not performed. The result is included.

The initial calibration run on 11/4/94 exceeds the acceptable QC limits for diethyl phthalate (31.079%). Since no samples were quantitated by this initial calibration, the data are not affected.

The continuing calibration run on 2/8/95 exceeds the acceptable QC limits for hexachlorocyclopentadiene (25.25%). Since this analyte was not detected in any of the samples, the data are not affected.

The continuing calibration run on 2/24/95 exceeds the acceptable QC limits for 1,2-dichlorobenzene (26.28%), 4-methyl phenol (31.58%), hexachloroethane (31.48%), di-n-octyl phthalate (31.43%) and 4-nitroaniline (28.77%). All detected values for sample 00914 should be regarded as estimated.

The continuing calibration run on 2/27/95 exceeds the acceptable QC limits for hexachloroethane (25.51%), hexachlorobutadiene (25.86%), fluorene (29.24%) and benzo (k) fluoranthene (33.55%). Since these analytes were not detected in any of the samples, the data are not affected.

The surrogates were not recovered for samples 00912 and 00913 due to the large dilutions necessary for the analysis. The data are not affected.

The continuing calibration run on 2/10/95 exceeded the acceptable QC limit for methylene chloride (32.44%), 1,2-dichloroethane-d₄, (39.23), 1,2 dichloroethane (30.59%), 4-methyl-2-pentanone (29.10%) and 2-hexanone (31.82%). Since no samples were quantitated by this calibration, the data are not affected.

One surrogate percent recovery exceeds the acceptable QC limits for blank G 6677. The data are not affected.

BNA TCLP Package E 78 (Drum Solids)

The initial calibration run on 2/7/95 on instrument F exceeds the acceptable QC limits for N-nitrosodimethyl amine (34.35%) and benzidine (65.36%). N-nitrosodimethyl amine was not detected in any of the associated samples and benzidine is no longer a CLP compound. The data are not affected.

The initial calibration run on 2/6/95 on instrument B exceeds the acceptable QC limits for benzoic acid (31.98%) and benzidine (60.96%). The data for benzoic acid should be regarded as estimated; benzidine is no longer a CLP compound.

The continuing calibration check standard run on 2/9/95 on instrument F exceeds the acceptable QC limits for 2,4,6-tribromophenol (45.20%), pyrene (28.12%), butyl benzyl phthalate (52.15%) and bis (2-ethylhexyl) phthalate (45.52%). The positive results and non-detected values for bis (2-ethylhexyl) phthalate should be regarded as estimated. The non detected values for butyl benzyl phthalate should also be regarded as estimated. The data for 2,4,6-tribromophenol and pyrene are not affected.

The continuing calibration check standard run on 2/11/95 on instrument F failed. Pentachlorophenol, a calibration check compound, exceeded the acceptable QC limit (36.48%). Since this standard was run for the blank and QA/QC samples only, an exception will be made. Positive results will not be qualified; however, undetected values reported in the blank will be considered as estimated for the following compounds because the percent difference was greater than 50: benzoic acid (70.43%), 2,4-dinitrophenol (52.01%), 4,6-dinitro-2-methyl phenol (59.87%) and butyl benzyl phthalate (60.09%).

The continuing calibration check standard run on 2/14/95 on instrument B was used to quantitate sample 00911 (DL) and 00916 (DL) for benzyl alcohol and benzoic acid. The percent differences for these analytes exceed the QC limits, 28.59% and 32.35%, respectively. The positive results for these compounds should be regarded as estimated only if the dilution values are used.

The result for benzoic acid in sample 00916 should be regarded as estimated because the concentration found in the diluted sample exceeds the highest linear standard.

The internal standard, perylene- d_{12} , is less than 50% of the area for perylene- d_{12} in the associated continuing calibration check standard. Because the sample is a QA sample, the data are not affected.

The surrogate, 2-fluorobiphenyl was either not found or below the acceptable QC limit in samples 00911, 00916, 00915 MS and 00915 MSD due to coeluting interferences. The associated data are not affected.

The matrix spike percent recovery for 1,2,4-trichlorobenzene exceeds thie acceptable QC limit.

The relative percent differences for 1,4-dichlorobenzene, 1,2,4-trichlorobenzene, 4-nitrophenol and 2,4-dinitrotoluene in the MS/MSD analysis exceeded the acceptable QC limits. The associated sample data are not affected.

Pesticide/PCB TCLP Package E 78 (Drum Solids)

The associated data are not affected.

The holding time was exceeded for all samples, however, as the holding time criterion does not apply to drum solids, the data are not affected.

Because the samples had been diluted, the surrogate, dibutyl chlorendate, was not recovered for samples 00911 10x and 00915 10x. Since these samples were used for confirmation purposes, the data are not affected.

The percent recoveries for lindane exceeded the acceptable QC limits for samples 00911 MS and 00911 MSD. The percent recoveries for heptachlor and endrin exceeded the acceptable QC limits for sample 00911 MS. The percent recoveries for dieldrin and 4,4'-DDT exceeded the acceptable QC limits for sample 00911 MSD. The data are not affected.

The MS/MSD results are not usable because of the high relative percent differences (44-124) obtained in the analysis.

Metals Package E 56

Mercury and matrix spike/matrix spike duplicate samples were not analyzed due to insufficient sample size.

At the time of collection, preservative was not added to the samples to be analyzed for metals.

SUMMARY of ABBREVIATIONS

В	The analyte was found i	n the blank	
BFB	Bromofluorobenzene		
BPQL	Below the Practical Qua	intitation Limit	
С	Centigrade		
D	(Surrogate Table) this v	alue is from a di	luted sample and was not calculated
	(Result Table) this result	lt was obtained f	rom a diluted sample
CLP	Contract Laboratory Pro	otocol	
COC	Chain of Custody		
CONC	Concentration		
CRDL	Contract Required Dete	ction Limit	
DFTPP	Decafluorotriphenylphosp	ohine	
DL	Detection Limit		
E	The value is greater than	n the highest line	ear standard and is estimated
EMPC	Estimated maximum pos	ssible concentration	on
J	The value is below the r	method detection	limit and is estimated
HHL	High Hazard Laboratory	, Brunswick, GA	
IDL	Instrument Detection L:	imit	
ISTD	Internal Standard		
MDL	Method Detection Limit	t	
MQL	Method Quantitation Li	imit	
MI	Matrix Interference		
MS	Matrix spike		
MSD	Matrix spike duplicate		
MW	Molecular weight		
NA	either Not Applicable of	r Not Available	
NC	Not Calculated		
ND	Not Detected		
NR	Not Requested		
NS	Not Spiked		
% D	Percent difference		
% REC	Percent Recovery		
PQL	Practical quantitation lin	mit	
PPBV	Parts per billion by volu	me	
QL	Quantitation Limit		
RPD	Relative percent differer	nce	
RSD	Relative Standard Devia	tion	
SIM	Selected Ion Mode		
m³	cubic meter	kg	kilogram
l(L)	liter	g	gram
dl	deciliter	cg	centigram
ml	milliliter	mg	milligram
ul	microliter	ug	microgram
		ng	nanogram

denotes a value that exceeds the acceptable QC limit

pg

Abbreviations that are specific to a particular table are explained in footnotes on that table

picogram

Revision 11/17/94

098\DEL\AR\9503\RE F ROC

48 🦸 🐧

Analytical Procedure for VOA in Water and Drum Liquids

A modified 524.2 method was used for the analysis of Volatile Organic Compounds in water and drum liquids. Samples were purged, trapped, and desorbed to a GC/MS system. Prior to purging, the samples were spiked with a three component surrogate mixture consisting of toluene-d₈, 4-bromofluorobenzene and 1,2-dichloroethane-d₄ and a three component internal standard mixture consisting of bromochloromethane, 1,4-difluorobenzene, and chlorobenzene-d₅. The following conditions and parameters were utilized:

1) Purge and Trap Unit: A Tekmar concentrator (LSC 2000) equipped with an autosampler (ALS2016) was utilized.

Purge and Trap parameters:

Purge

10 min at 25° C

Dry Purge

2 min at 25° C

Desorb
Purge Flow Rate

4 min at 250° C

40 ml/min

Desorb Preheat 245° C Bake 8 min

8 min at 260° C

Trap - VOCARB 4000 (Supelco) which consists of four adsorbent beds: Carbopack B (graphitized carbon 60/80 mesh), Carbopack C (graphitized carbon 60/80 mesh), Carboxen-1000 (60/80 mesh), and Carboxen-1001 (60/80 mesh).

2) GC/MS System: A Hewlett Packard 5970 GC/MSD equipped with an RTE-A data system was used.

GC/MS parameters:

GC Column: thickness.

30 meter x 0.53mm ID, RTx-Volatiles (Restek Corp.) column with 3.0 um

GC Temperature

5 min at 10° C;

6° C/min to 140° C

Program

0.1 min at 140° C;

12° C/min to 160° C

5 min at 160° C

GC Flow Rate

Helium at 10 ml/min

GC/MS Interface - Glass jet separator with 30 ml/min helium make-up gas at 250° C.

Mass Spectrometer - Electron Impact Ionization at a nominal electron energy of 70 electron volts, scanning from 35-300 amu at one scan/sec.

Computer - Preprogrammed to plot Extracted Ion Current Profile (EICP); capable of integrating ions and plotting abundances vs time or scan number. A library search (NBS-Wiley) for tentatively identified compounds was performed on samples.

The GC/MS system was calibrated using 6 VOA standards at 5, 20, 50, 100, 150, and 200 ug/l. Before analysis each day, the system was tuned with 50 ng BFB and passed a continuing calibration check when analyzing a 50 ug/l standard mixture in which the responses were evaluated by comparison to the average response of the calibration curve.

The results for water are listed in Table 1.1 and the Tentatively Identified Compounds for water are listed in Table 1.2. The results for drum liquids are listed in Table 1.6 and the Tentatively Identified

The concentrations of the analytes were calculated using the following equation:

$$C_{u} = \frac{A_{x} \times I_{x}}{A_{x} \times RF \text{ (or } RF_{ave}) \times V_{o}}$$

Compounds for drum liquids are listed in Table 1.7.

where

 C_u = concentration of the target analyte (ug/L)

 A_x = Area of the target analyte

 I_{is} = ng of specific internal standard

 A_{is} = Area of the specific internal standard

RF = Response Factor

RF_{ave} = average Response Factor

V_o = volume of sample purged (ml), taking into account dilutions

The average Response Factor is used when a sample is associated with an initial calibration curve. The Response Factor is used when a sample is associated with a continuing calibration curve.

Response Factor calculation:

The response factor (RF) for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_c \times I_u}{A_u \times I_c}$$

where,

RF = Response factor for a specific analyte

 A_c = Area of the analyte in the standard

I₁₅ = Mass of the specific internal standard

A_s = Area of the specific internal standard

 I_c = Mass of the analyte in the standard

$$RF_{ave} = \frac{RF_1 + ... + RF_n}{n}$$

where

n = number of Samples

Revision of 6/28/94

2 1 32/3

Analytical Procedure for VOA in Soil

A modified 524.2 method was used for the analysis of Volatile Organic Compounds in soil. Samples were purged, trapped, and desorbed to a GC/MS system. Prior to purging, the samples were spiked with a three component surrogate mixture consisting of toluene-d₈, 4-bromofluorobenzene and 1,2-dichloroethane-d₄ and a three component internal standard mixture consisting of bromochloromethane, 1,4-diffuorobenzene, and chlorobenzene-d₅. The following conditions and parameters were utilized:

1) Purge and Trap Unit: A Tekmar concentrator (LSC 2000) equipped with an autosampler (ALS2016) was utilized.

Purge and Trap parameters:

Purge 10 min at 25° C Dry Purge 2 min at 25° C
Desorb 4 min at 250° C Desorb Preheat 245° C
Purge Flow Rate 40 ml/min Bake 8 min at 260° C

Trap - VOCARB 4000 (Supelco) which consists of four adsorbent beds: Carbopack B (graphitized carbon 60/80 mesh), Carbopack C (graphitized carbon 60/80 mesh), Carboxen-1000 (60/80 mesh), and Carboxen-1001 (60/80 mesh).

2) GC/MS System: A Hewlett Packard 5970 GC/MSD equipped with an RTE-A data system was used.

GC Column: 30 meter x 0.53mm ID, RTx - Volatiles (Restek Corp.) column with 3um thickness.

GC Temperature Program:

5 min at 10° C;

6° C/min to 140° C

0.1 min at 140° C;

12° C/min to 160° C

5 min at 160° C

GC Flow Rate:

Helium at 10 ml/min

GC/MS Interface:

Glass jet separator with 30 ml make-up gas at 250° C.

Mass Spectrometer: Electron Impact Ionization at a nominal electron energy of 70 electron volts, scanning from 35-300 amu at one scan/sec.

Computer - Preprogrammed to plot Extracted Ion Current Profile (EICP); capable of integrating ions and plotting abundances vs time or scan number. A library search (NBS-Wiley) for tentatively identified compounds was performed on samples.

The GC/MS system was calibrated using 6 VOA standards at 5, 20, 50, 100, 150, and 200 ug/l. Before analysis each day, the system was tuned with 50 ng BFB and passed a continuing calibration check when analyzing a 50 ug/l standard mixture in which the responses were evaluated by comparison to the average response of the calibration curve.

The medium level soil extracts were analyzed by extracting 5.0 g soil with 5 ml methanol, diluting an aliquot with 5 ml water and analyzing the solution by the purge and trap method. The results are in Table 1.3; the tentatively identified compounds are listed in Table 1.4. The concentrations of the analytes were calculated using the following equation:

$$C_{u} = \frac{DF \times A_{x} \times I_{u}}{A_{u} \times RF \text{ (or } RF_{ave}) \times W_{s} \times D}$$

where

C_u = Concentration of target analyte (ug/kg) on a dry weight basis

DF = Dilution Factor

 A_x = Area of the target analyte

I_s = mass of specific internal standard (ng) A_s = Area of the specific internal standard

RF = Response Factor

RF_{ave} = average Response Factor W, = Weight of sample (g) D = Decimal percent solids

The average Response Factor is used when a sample is associated with an initial calibration curve. The Response Factor is used when a sample is associated with a continuing calibration curve.

Response Factor calculation:

The response factor (RF) for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_c \times I_s}{A_s \times I_c}$$

where.

RF = Response factor for a specific analyte A_c = Area of the analyte in the standard I_s = Mass of the specific internal standard A_u = Area of the specific internal standard I_c = Mass of the analyte in the standard

$$RF_{ave} = \frac{RF_1 + ... + RF_n}{n}$$

and

n = number of Samples

Revision of 6/28/94

2.2 - 12.0

Analytical Procedure for TCLP VOA

The subcontract laboratory determined the concentration of VOA compounds in the TCLP leachate by preparing the samples according to Method 1311 and analyzing the leachates according to Method 8240A. Both methods are given in "Test Methods for Evaluating Solid Wastes, Physical and Chemical Methods: 3rd Edition, USEPA SW846, 1986, Final Update I, July 1992". The results of the analysis for the sediment samples are listed in Table 1.5. The results of the analysis for the drum solid samples are listed in Table 1.8.

Analytical Procedure for BNA in Water, TCLP Extracts and Drum Liquids

Extraction Procedure

Prior to extraction, each sample was spiked with a six component surrogate mixture consisting of nitrobenzene-d₅, 2-fluorobiphenyl, terphenyl-d₁₄, phenol-d₅, 2-fluorophenol, and 2,4,6-tribromophenol. One liter of sample was extracted according to Method 625, Section 10, as outlined in the Federal Register Vol. 49, #209, Friday, Oct. 26, 1984. The extracts were combined, concentrated to 1.0 ml, an internal standards mixture consisting of 1,4-dichlorobenzene-d₄, naphthalene-d₈, acenaphthene-d₁₀, phenanthrene-d₁₀, chrysene-d₁₂, and perylene-d₁₂ was added, and analyzed.

Analytical Procedure

An HP 5995C Gas Chromatograph/Mass Spectrometer (GC/MS), equipped with a 7673A autosampler and controlled by an HP-1000 RTE-6/VM computer was used to analyze the samples.

The instrument conditions were:

Column	Restek Rtx-5 (crossbonded SE-54)
	30 meter x 0.32mm ID, 0.50 μm
	film thickness
Injection Temperature	290° C
Transfer Temperature	290° C
Source Temperature	240° C
Analyzer Temperature	240° C
Temperature Program	40°C for 3 min
•	8° C/min to 295° C
	hold for 12 min
Splitless Injection	Split time $= 1.00 \text{min}$
Injection Volume	1 ul

The GC/MS system was calibrated using 5 BNA standards at 20, 50, 80, 120, and 160 μ g/ml. Before analysis each day, the system was tuned with 50 ng decafluorotriphenylphosphine (DFTPP) and passed a continuing calibration check when analysing a 50 μ g/ml standard mixture in which the responses were evaluated by comparison to the average response of the calibration curve.

The BNA results for the water samples are listed in Table 1.9; the tentatively identified compounds are listed in Table 1.10. The BNA results for the TCLP extracts are listed in Table 1.13; the tentatively identified compounds are listed in Table 1.14. The BNA results for the drum liquids are listed in Table 1.15; the tentatively identified compounds are listed in Table 1.16. The concentration of the detected compounds was calculated using the following equation:

$$C_{u} = \frac{DFxA_{u}xI_{is}xV_{t}}{A_{is}xRF(orRF_{ane})xV_{i}xV_{o}}$$

where

 C_u = Concentration of target analyte ($\mu g/L$)

DF = Dilution Factor

 A_u = Area of target analyte

I_s = Mass of specific internal standard (ng)

 V_i = Volume of extract (μl)

A_s = Area of specific internal standard

RF = Response Factor (unitless) RF₁₀₀ = average Response Factor

 V_i = Volume of extract injected (μl)

 V_{o} = Volume of sample (ml)

The RF_{ave} is used when a sample is associated with an initial calibration curve. The RF is used when a sample is associated with a continuing calibration curve.

Response Factor calculation:

The RF for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_c \times I_{is}}{A_{is} \times I_c}$$

where

RF = Response factor for a specific analyte A_c = Area of the analyte in the standard I_s = Mass of the specific internal standard A_s = Area of the specific internal standard I_c = Mass of the analyte in the standard

$$RF_{ave} = \frac{RF_1 + ... + RF_n}{n}$$

and

n = number of Samples

Rev. 7/11/94

Analytical Procedure for BNA in Soil

2 1 3233

Extraction Procedure

Prior to extraction each sample was spiked with a six component surrogate mixture consisting of nitrobenzene- d_5 , 2-fluorobiphenyl, terphenyl- d_{14} , phenol- d_5 , 2-fluorophenol, and 2,4,6-tribromophenol. Thirty grams of sample was mixed with 30 g anhydrous sodium sulfate, and Soxhlet extracted for 16 hours with 300 ml of 1:1 acetone:methylene chloride. The extract was concentrated to 1.0 ml., an internal standard mixture consisting of 1,4-dichlorobenzene- d_4 , naphthalene- d_8 , acenaphthene- d_{10} , phenanthrene- d_{10} , chrysene- d_{12} , and perylene- d_{12} was added, and analyzed.

Analysis Procedure

An HP 5995C Gas Chromatograph/Mass Spectrometer (GC/MS), equipped with a 7673A autosampler and controlled by an HP-1000 RTE-6/VM computer was used to analyze the samples.

The instrument conditions were:

Column	Restek Rtx-5 (crossbonded SE-54) 30 meter x 0.32mm ID, 0.50 µm film thickness
Injection Temperature	290° C
Transfer Temperature	290° C
Source Temperature	240° C
Analyzer Temperature	240° C
Temperature Program	40°C for 3 min
	8° C/min to 295° C
	hold for 12 min
Splitless Injection	Split time $= 1.00 \text{min}$
Injection Volume	1 μl

The GC/MS system was calibrated using 5 BNA standard mixtures at 20, 50, 80, 120, and 160 μ g/ml. Before analysis each day, the system was tuned with 50 ng decafluorotriphenylphosphine (DFTPP) passed a continuing calibration check when analyzing a 50 μ g/ml standard mixture in which the responses were evaluated by comparison to the average response of the calibration curve.

The BNA results, based on dry weight, are listed in Table 1.11; the tentatively identified compounds are listed in Table 1.12. The concentration of the detected compounds was calculated using the following equation:

$$C_{u} = \frac{DFxA_{u}xI_{ix}xV_{t}}{A_{ix}xRF(orRF_{ave})xV_{i}xWxD}$$

where

 C_{μ} = Concentration of target analyte ($\mu g/Kg$)

DF = Dilution Factor

A_u = Area of target analyte

 I_{is} = Mass of specific internal standard (ng)

 V_i = Volume of extract (μ l)

A_{is} = Area of specific internal standard

RF = Response Factor (unitless) RF_{ave} = average Response Factor

 V_1 = Volume of extract injected (μ l)

W = Weight of sample (g)
D = Decimal per cent solids

The RF_{ave} is used when a sample is associated with an initial calibration curve. The RF is used when a sample is associated with a continuing calibration.

Response Factor calculation:

The RF for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_c \times I_{is}}{A_{is} \times I_c}$$

where

RF = Response factor for a specific analyte

A_c = Area of the analyte in the standard

I_s = Mass of the specific internal standard

A_s = Area of the specific internal standard

I_c = Mass of the analyte in the standard

$$RF_{ave} = \frac{RF_1 + \dots + RF_n}{n}$$

and

n = number of Samples Revision of 7/08/94

Analytical Procedure for TCLP BNA

The subcontract laboratory determined the concentration of BNA compounds in the TCLP leachate by preparing the samples according to Method 1311 and analyzing the leachates according to Method 8270. Both methods are given in "Test Methods for Evaluating Solid Wastes, Physical and Chemical Methods: 3rd Edition, USEPA SW846, 1986, Final Update I, July 1992". The results of the analysis for the drum solid samples are listed in Table 1.17.

Analytical Procedure for Pesticides and PCBs in Water

Extraction Procedure

One liter of sample was spiked with a surrogate solution consisting of tetrachloro-m-xylene and decachlorobiphenyl, and was extracted three times with 60 ml portions of methylene chloride. The combined extracts were filtered, concentrated to 10 ml, solvent exchanged with 60 ml hexane, and the hexane concentrated to 1.0 ml.

Gas Chromatographic Analysis

The extract was analyzed for pesticides using simultaneous dual column injections. The analysis was done on an HP 5890 GC/ECD system, equipped with an HP 7673A automatic sampler, and controlled with an HP-ChemStation. The following conditions were employed:

First Column DB-608, 30 meter, 0.53mm fused silica

capillary, 0.83 µm film thickness

Injector Temperature 250° C
Detector Temperature 325° C

Temperature Program 150°C for 1 minute 7°C/min to 265°C

18 min at 265°

Second Column Rtx-1701, 30 meter, 0.53mm fused silica

capillary, 0.50 µm film thickness

Injector Temperature 250° C
Detector Temperature 325° C

Temperature Program 150° C for 1 minute 7°C/min to 265°C

18 min at 265°

The gas chromatographs were calibrated using 5 pesticide standards at 20, 50, 100, 200, and 500 μ g/L. The response from each mixture were used to calculate the response factors (RF) of each analyte. The average RF was used to calculate the concentrations of the pesticides in the samples. Quantification was based on the DB-608 column (signal 1), and identity of the analyte was confirmed using the Rtx-1701 column (signal 2). A fingerprint gas chromatogram was run using each of the seven Aroclor mixtures, and toxaphene; calibration curves were run only if a particular Aroclor, or toxaphene was found in the sample.

The Pesticide/PCB results, listed in Table 1.18, were calculated from the following formula:

$$C_{ii} = \frac{DFxA_{ii}xV_{i}}{RF_{ane}xV_{i}xV_{i}}$$

where

 C_u = Concentration of analyte ($\mu g/L$)

DF = Dilution Factor A_u = Area or peak height V_i = Volume of sample (ml) RF_{ave} = Average response factor V_i = Volume of extract injected (μ l)

 $V_s = Sample volume (ml)$

Response Factor calculation:

The RF for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_n}{total \ pg \ injected}$$

where

A_u = Area or peak height

and

$$RF_{ave} = \frac{RF_1 + ... + RF_n}{n}$$

where

n = number of samples

Revision 7/11/94

2 1 3213

Analytical Procedure for Pesticides and PCBs in Soil

Extraction Procedure

Thirty grams of sample was spiked with a surrogate solution consisting of tetrachloro-m-xylene and decachlorobiphenyl, 30 g anhydrous sodium sulfate and Soxhlet extracted with 16 hours with 300 ml 1:1 hexane: acetone. The extract was concentrated to 5.0 ml.

Gas Chromatographic Analysis

The extract was analyzed for pesticides and PCBs using simultaneous dual column injections. The analysis was done on an HP 5890 GC/ECD system, equipped with an HP 7673A automatic sampler, and controlled with an HP-CHEM STATION. The following conditions were employed:

First Column DB-608, 30 meter, 0.53mm fused silica

capillary, 0.83 µm film thickness

Injector Temperature 250° C
Detector Temperature 325° C

Temperature Program 150°C for 1 minute

7°C/min to 265°C 18 min at 265°

Second Column Rtx-1701, 30 meter, 0.53mm fused silica

capillary, 0.50 µm film thickness

Injector Temperature 250° C
Detector Temperature 325° C

Temperature Program 150° C for 1 minute 17°C/min to 265°C

18 min at 265°

The gas chromatographs were calibrated using 5 pesticide standards at 20, 50, 100, 200, and 500 μ g/L. The results from each mixture were used to calculate the response factor (RF) of each analyte and the average Response Factor was used to calculate the concentration of pesticide in the sample. Quantification was based on the DB-608 column (signal 1) and the identity of the analyte was confirmed using the Rtx-1701 column (signal 2). A fingerprint chromatogram was run using each of the seven Aroclor mixtures and toxaphene; calibration curves were run only if a particular Aroclor or toxaphene was found in the sample.

The pesticide/PCB results, listed in Table 1.19, are calculated by using the following formula:

$$C_u = \frac{DFxA_uxV_t}{RF_{out}xV_txWxD}$$

where

 C_u = Concentration of analyte (mg/Kg)

DF = Dilution Factor A_u = Area or peak height V_t = Volume of sample (ml) RF_{ave} = Average response factor V_i = Volume of extract injected (μ l)

W = Weight of sample (g)
D = Decimal per cent solids

Response Factor calculation:

The RF for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_u}{total \ pg \ injected}$$

where

A_u = Area or peak height

and

$$RF_{ave} = \frac{RF_1 + \dots + RF_n}{n}$$

where

n = number of samples

Revision 7/11/94

2 1 3210

Analytical Procedure for Pesticides and PCBs in Drum Liquids

Extraction Procedure

One hundred and eighty ml of sample was spiked with a surrogate solution consisting of tetrachloro-m-xylene and decachlorobiphenyl, and was extracted three times with 60 ml portions of methylene chloride. The combined extracts were filtered, concentrated to 10 ml, solvent exchanged with 60 ml hexane, and the hexane concentrated to 1.0 ml.

Gas Chromatographic Analysis

The extract was analyzed for pesticides using simultaneous dual column injections. The analysis was done on an HP 5890 GC/ECD system, equipped with an HP 7673A automatic sampler, and controlled with an HP-ChemStation. The following conditions were employed:

First Column	DB-608, 30 meter, 0.53mm fused silica
	agaillan, 0.92 um film thiolmaga

capillary, 0.83 µm film thickness

Injector Temperature 250° C
Detector Temperature 325° C

Temperature Program 150°C for 1 minute 7°C/min to 265°C 18 min at 265°

Second Column Rtx-1701, 30 meter, 0.53mm fused silica

capillary, 0.50 µm film thickness

Injector Temperature 250° C Detector Temperature 325° C

Temperature Program 150° C for 1 minute 7°C/min to 265°C

18 min at 265°

The gas chromatographs were calibrated using 5 pesticide standards at 20, 50, 100, 200, and 500 $\mu g/L$. The response from each mixture were used to calculate the response factors (RF) of each analyte. The average RF was used to calculate the concentrations of the pesticides in the samples. Quantification was based on the DB-608 column (signal 1), and identity of the analyte was confirmed using the Rtx-1701 column (signal 2). A fingerprint gas chromatogram was run using each of the seven Aroclor mixtures, and toxaphene; calibration curves were run only if a particular Aroclor, or toxaphene was found in the sample.

The Pesticide/PCB results, listed in Table 1.21, were calculated from the following formula:

$$C_{u} = \frac{DFxA_{u}xV_{t}}{RF_{ave}xV_{t}xV_{s}}$$

where

 C_u = Concentration of analyte $(\mu g/L)$

DF = Dilution Factor

A_u = Area or peak height

V_t = Volume of sample (ml)

RF_{ave} = Average response factor

 V_i = Volume of extract injected (μ l)

 $V_s = Sample volume (ml)$

Response Factor calculation:

The RF for each specific analyte is quantitated based on the area response from the continuing calibration check as follows:

$$RF = \frac{A_u}{total \ pg \ injected}$$

where

A_u = Area or peak height

and

$$RF_{ave} = \frac{RF_1 + ... + RF_n}{n}$$

where

n = number of samples

Revision 7/11/94

2 : 1212

Analytical Procedure for TCLP Pesticide/PCBs

The subcontract laboratory determined the concentration of pesticides and PCBs in the TCLP leachate by preparing the samples according to Method 1311 and analyzing the leachates according to Method 8080. Both methods are given in "Test Methods for Evaluating Solid Wastes, Physical and Chemical Methods: 3rd Edition, USEPA SW846, 1986, Final Update I, July 1992". The results of the analysis for the drum solid samples are listed in Table 1.22.

Analytical Procedure for Metals in Water

A 45 ml aliquot of sample was mixed with 5.0 ml concentrated nitric acid, placed in an acid rinsed Teflon container, capped with a Teflon lined cap, and digested according to SW-846, Method 3015, in a CEM MDS-2100 microwave oven (the first stage was at 99% power for 10 minutes and the second stage was at 70% power for 10 minutes). After digestion, the sample was analysed for all metals, except mercury, by USEPA SW-846 Method 7000/6010.

Mercury was analyzed separately on a Varian SpectrAA-300 Atomic Absorption Spectrophotometer equipped with a Varian VGA-76 vapor gas analyser using modified method 7470 as given by "Test Method for Evaluating Solid Waste, Sept. 1986," USEPA SW-846. Results of the analyses of the water samples are listed in Table 1.23.

Analytical Procedure for Metals in Soil

One gram of sample, weighed to 0.01 g accuracy, was thoroughly mixed with 10 ml 1:1 nitric acid:water, and digested according to SW-846, Method #3050 and analyzed according to Method 7000/6010.

Mercury was analyzed separately on a Varian SpectrAA-20 Atomic Absorption Spectrophotometer equipped with a Varian VGA-76 vapor gas analyzer using method 7471 as given by "Test Methods for Evaluating Solid Waste", USEPA, SW-846, September, 1986. Results of the analyses are listed in Table 1.24.

Analytical Procedure for TCLP Metals

A 45 ml aliquot of sample was mixed with 5.0 ml concentrated nitric acid, placed in an acid rinsed Teflon container, capped with a Teflon lined cap, and digested according to SW-846, Method 3015, in a CEM MDS-2100 microwave oven (the first stage was at 99% power for 10 minutes and the second stage was at 70% power for 10 minutes). After digestion, the sample was analysed for all metals, except mercury, by USEPA SW-846 Method 7000/6010.

Mercury was analyzed separately on a Varian SpectrAA-300 Atomic Absorption Spectrophotometer equipped with a Varian VGA-76 vapor gas analyser using modified method 7470 as given by "Test Method for Evaluating Solid Waste, Sept. 1986," USEPA SW-846. Results of the analyses of the TCLP leachates are listed in Table 1.25.

Analytical Procedure for Metals in Drum Liquid Samples

A 25 ml aliquot of sample was allowed to evaporate to dryness on a hot plate, mixed with 5.0 ml concentrated nitric acid, placed in an acid rinsed Teflon container, capped with a Teflon lined cap, and digested according to SW-846, Method 3010. After digestion, the sample was diluted to 50 ml with ASTM Type II water and analysed for all metals by USEPA SW-846 procedures. The metals analyses were done either on a Varian SpectrAA-400Z Atomic Absorption Spectrophotometer or on an Inductively Coupled Argon Plasma (ICAP) emission spectrometer. Results of the analyses of the samples are listed in Table 1.26.

Analytical Procedure for TCLP Metals in Drum Solids

The subcontract laboratory determined the concentration of metals in the TCLP leachate by preparing the samples according to Method 1311 and analyzing the leachates according to Method 7000. Both methods are given in "Test Methods for Evaluating Solid Wastes, Physical and Chemical Methods: 3rd Edition, USEPA SW846, 1986, Final Update I, July 1992". The results of the analysis for the drum solid samples are listed in Table 1.27.

Analytical Procedure for BTU

The subcontract laboratory determined the BTU value of the samples using ASTM Method D-240. The results of the analysis are listed in Table 1.28.

Analytical Procedure for Ash

The subcontract laboratory determined the ash content of the samples using ASTM Method D-482. The results of the analysis are listed in Table 1.28.

Analytical Procedure for Chlorine

The subcontract laboratory determined the chlorine content of the samples using ASTM Method D-808. The results of the analysis are listed in Table 1.28.

Analytical Procedure for Sulfur

The subcontract laboratory determined the sulfur content of the samples using ASTM Method D-129. The results of the analysis are listed in Table 1.28.

Analytical Procedure for Moisture

The subcontract laboratory determined the moisture content of the samples using ASTM Method D-95. The results of the analysis are listed in Table 1.28.

2 0 3203

Table 1.1 Results of the VOA Analysis for the Water Samples WA # 0-098 Pier Drum

SAMPLE # : LOCATION : COLLECTED : ANALYZED : INJECTED : FILE # : DIL. FACT : UNIT :	01/13/9: 14:40 ^V2426 1 ug/L		00645 TRIP BL/ 01/11/99 01/13/99 15:40 ^V2427 1 ug/L	5	00638 N PIT 01/11/9 01/13/9 17:42 ^V2430 1 ug/L	
COMPOUND	CONC.	MDL	CONC.	MDL	CONC.	MDL
Dichlorodifluoromethane	ND	1.0	ND	1.0	ND	1.0
Chloromethane	ND	1.0	ND	1.0	ND	1.0
Vinyl Chloride	ND	1.0	ND	1.0	ND	1.0
Bromomethane	ND	2.0	ND	2.0	ND	2.0
Chloroethane	ND	1.0	ND	1.0	ND	1.0
Trichlorofluoromethane	ND	1.0	ND	1.0	ND	1.0
Acetone	ND	2.0	ND	2.0	ND	2.0
1,1-Dichloroethene	ND	1.0	ND	1.0	ND	1.0
Carbon Disulfide	ND	1.0	ND	1.0	ND	1.0
Methylene Chloride	ND	1.0	ND	1.0	ND	1.0
Methyl-tertiary-butylether	ND	1.0	ND	1.0	ND	1.0
trans-1,2-Dichloroethene	ND	1.0	ND	1.0	ND	1.0
1,1-Dichloroethane	ND	1.0	ND	1.0	ND	1.0
2-Butanone	ND	4.0 1.0	ND	4.0	ND	4.0
2,2-Dichloropropane	ND ND	1.0	ND	1.0	ND	1.0
cis-1,2-Dichloroethene			ND	1.0	ND	1.0
Chloroform	ND	1.0	ND	1.0	ND	1.0
1,1-Dichloropropene	ND	1.0	ND	1.0	ND	1.0
1,2-Dichloroethane	ND ND	1.0 1.0	ND ND	1.0	ND	1.0
1,1,1-Trichloroethane Carbon Tetrachloride	ND ND	1.0	ND	1.0 1.0	ND	1.0
Benzene	ND	1.0	ND	1.0	ND ND	1.0 1.0
Trichloroethene	ND	1.0	ND	1.0	ND	1.0
1.2-Dichloropropane	ND	1.0	ND	1.0	ND	1.0
Dibromomethane	ND	1.0	ND	1.0	ND	1.0
Bromodichloromethane	ND	1.0	ND	1.0	ND	1.0
cis-1,3-Dichloropropene	ND	1.0	ND	1.0	ND	1.0
trans-1,3-Dichloropropene	ND	1.0	ND	1.0	ND	1.0
1,1,2-Trichloroethane	ND	1.0	ND	1.0	ND	1.0
1,3-Dichloropropane	ND	1.0	ND	1.0	ND	1.0
Dibromochloromethane	ND	1.0	ND	1.0	ND	1.0
1,2-Dibromoethane	ND	1.0	ND	1.0	ND	1.0
Bromoform	ND	1.0	ND	1.0	ND	1.0
4-Methyl-2-Pentanone	ND	2.0	ND	2.0	ND	2.0
	• • • • • • • • • • • • • • • • • • • •			• • • • • •		

2 (2) 3297

Table 1.1 (Cont) Results of the VOA Analysis for the Water Samples WA # 0-098 Pier Drum

	SAMPLE # : LOCATION : COLLECTED : ANALYZED : INJECTED : FILE : DIL. FACT :	01/13/9 14:40 ^V2426		00645 TRIP BL 01/11/9 01/13/9 15:40 ^V2427	5	00638 N PIT 01/11/9 01/13/9 17:42 ^V2430 1	
	UNIT :	ug/L		ug/L		ug/L	
COMPOUND		CONC.	MDL	CONC.	MDL	CONC.	MDL
Toluene		ND	1.0	ND	1.0	ND	1.0
2-Hexanone		ND	2.0	ND	2.0	ND	2.0
Tetrachloroethene		ND	1.0	ND	1.0	ND	1.0
Chlorobenzene		ND	1.0	ND	1.0	ND	1.0
1,1,1,2-Tetrachlor	oethane	ND	1.0	ND	1.0	ND	1.0
Ethylbenzene		ND	1.0	ND	1.0	ND	1.0
p & m-Xylene		ND	1.0	ND	1.0	ND	1.0
o-Xylene		ND	1.0	ND	1.0	ND	1.0
Styrene		ND	1.0	ND	1.0	ND	1.0
1sopropylbenzene		ND	1.0	ND	1.0	ND	1.0
1,1,2,2-Tetrachlor		ND	1.0	ND	1.0	ND	1.0
1,2,3-Trichloropro	pane	ND	1.0	ND	1.0	ND	1.0
Bromobenzene		ND	1.0	ND	1.0	ND	1.0
n-Propylbenzene		ND	1.0	ND	1.0	ND	1.0
2-Chlorotoluene		ND	1.0	ND	1.0	ND	1.0
4-Chlorotoluene		ND	1.0	ND	1.0	ND	1.0
1,3,5-Trimethylben	zene	ND	1.0	ND	1.0	ND	1.0
tert-Butylbenzene		ND	1.0	ND	1.0	ND	1.0
1,2,4-Trimethylben:	zene	ND	1.0	ND	1.0	ND	1.0
sec-Butylbenzene		ND	1.0	ND	1.0	ND	1.0
1,3-Dichlorobenzen	e	ND	1.0	ND	1.0	ND	1.0
p-Isopropyltoluene		ND	1.0	ND	1.0	ND	1.0
1,4-Dichlorobenzen	2	ND	1.0	ND	1.0	ND	1.0
1,2-Dichlorobenzen	P	ND	1.0	ND	1.0	ND	1.0
n-Butylbenzene		ND	1.0	ND	1.0	ND	1.0
1,2-Dibromo-3-Chlo		ND	1.0	ND	1.0	ND	1.0
1,2,4-Trichloroben	zene	ND	1.0	ND	1.0	ND	1.0
Naphthalene		ND	1.0	ND	1.0	ND	1.0
Hexachlorobutadien		ND	1.0	ND	1.0	ND	1.0
1,2,3-Trichloroben	zene	ND	1.0	ND	1.0	ND	1.0

Table 1.1 (Cont) Results of the VOA Analysis for the Water Samples WA # 0-098 Pier Drum

SAMPLE # :	LAB BLANK	00640
LOCATION :		FIELD BLANK
COLLECTED :		01/11/95
ANALYZED :	01/14/95	01/14/95
INJECTED :	12:49	14:11
FILE # :	^V2441	^v2443
DIL. FACT.:	1	1
UNIT :	ug/L	ug/L

UNII	: ug/L		ug/L	
COMPOUND	CONC.	MDL	CONC.	MDL
Dichlorodifluoromethane	ND	1.0	ND	1.0
Chloromethane	ND	1.0	ND	1.0
Vinyl Chloride	ND	1.0	ND	1.0
Bromomethane	ND	2.0	ND	2.0
Chloroethane	ND	1.0	ND	1.0
Trichlorofluoromethane	ND	1.0	ND	1.0
Acetone	ND	2.0	ND	2.0
1,1-Dichloroethene	ND	1.0	ND	1.0
Carbon Disulfide	ND	1.0	ND	1.0
Methylene Chloride	ND	1.0	ND	1.0
Methyl-tertiary-butylether	ND	1.0	ND	1.0
trans-1,2-Dichloroethene	ND	1.0	ND	1.0
1,1-Dichloroethane	ND	1.0	ND	1.0
2-Butanone	ND	4.0	ND	4.0
2,2-Dichloropropane	ND	1.0	ND	1.0
cis-1,2-Dichloroethene	ND	1.0	ND	1.0
Chloroform	ND	1.0	ND	1.0
1,1-Dichloropropene	ND	1.0	ND	1.0
1,2-Dichloroethane	ND	1.0	ND	1.0
1,1,1-Trichloroethane	ND	1.0	ND	1.0
Carbon Tetrachloride	ND	1.0	ND	1.0
Benzene	ND	1.0	ND	1.0
Trichloroethene	ND	1.0	ND	1.0
1,2-Dichloropropane	ND	1.0	ND	1.0
Dibromomethane	ND	1.0	ND	1.0
Bromodichloromethane	ND	1.0	ND	1.0
cis-1,3-Dichloropropene	ND	1.0	ND	1.0
trans-1,3-Dichloropropene	ND	1.0	ND	1.0
1,1,2-Trichloroethane	ND	1.0	ND	1.0
1,3-Dichloropropane	ND	1.0	ND	1.0
Dibromochloromethane	ND	1.0	ND	1.0
1,2-Dibromoethane	ND	1.0	ND	1.0
Bromoform	ND	1.0	ND	1.0
4-Methyl-2-Pentanone	ND	2.0	ND	2.0

Table 1.1 (Cont) Results of the VOA Analysis for the Water Samples WA # 0-098 Pier Drum

SAMPLE # :	LAB BLANK	00640
LOCATION :		FIELD BLANK
COLLECTED :		01/11/95
ANALYZED :	01/14/95	01/14/95
INJECTED :	12:49	14:11
FILE :	^V2441	^V2443
DIL. FACT.:	1	1
UNIT :	ug/L	ug/L

COMPOUND	CONC.	MDL	CONC.	MDL
Toluene	ND	1.0	ND	1.0
2-Hexanone	ND	2.0	ND	2.0
Tetrachloroethene	ND	1.0	ND	1.0
Chlorobenzene	ND	1.0	ND	1.0
1,1,1,2-Tetrachloroethane	ND	1.0	ND	1.0
Ethylbenzene	ND	1.0	ND	1.0
p & m-Xylene	ND	1.0	ND	1.0
o-Xylene	ND	1.0	ND	1.0
Styrene	ND	1.0	ND	1.0
I sopropy l benzene	ND	1.0	ND	1.0
1,1,2,2-Tetrachloroethane	ND	1.0	ND	1.0
1,2,3-Trichloropropane	ND	1.0	ND	1.0
Bromobenzene	ND	1.0	ND	1.0
n-Propylbenzene	ND	1.0	ND	1.0
2-Chlorotoluene	ND	1.0	ND	1.0
4-Chlorotoluene	ND	1.0	ND	1.0
1,3,5-Trimethylbenzene	ND	1.0	ND	1.0
tert-Butylbenzene	ND	1.0	ND	1.0
1,2,4-Trimethylbenzene	ND	1.0	ND	1.0
sec-Butylbenzene	ND	1.0	ND	1.0
1,3-Dichlorobenzene	ND	1.0	ND	1.0
p-Isopropyltoluene	ND	1.0	ND	1.0
1,4-Dichlorobenzene	ND	1.0	ND	1.0
1,2-Dichlorobenzene	ND	1.0	ND	1.0
n-Butylbenzene	ND	1.0	ND	1.0
1,2-Dibromo-3-Chloropropane	ND	1.0	ND	1.0
1,2,4-Trichlorobenzene	ND	1.0	ND	1.0
Naphthalene	ND	1.0	ND	1.0
Hexachlorobutadiene	ND	1.0	ND	1.0
1,2,3-Trichlorobenzene	ND	1.0	ND	1.0

2 1 9299

Table 1.1 (Cont) Results of the VOA Analysis for the Water Samples $$\rm WA~\#~0\text{-}098~Pier~Drum$$

	SAMPLE # : LOCATION : COLLECTED : ANALYZED : INJECTED : FILE # : DIL. FACT.: UNIT :	01/18/95 00:24 ^V2472 1 ug/L		00636 MW-S DUP 01/11/95 01/18/95 01:04 ^V2473 1 ug/L		00641 MW-N 01/11/95 01/18/95 01:45 ^V2474 1 ug/L		00635 MW-S 01/11/95 01/18/95 02:25 ^V2475 1 ug/L		00634 MW-W 01/11/95 01/18/95 03:46 ^V2477 1 ug/L	
COMPOUND		CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL
Dichlorodifluorome	thane	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Chloromethane		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Vinyl Chloride		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Bromomethane		ND	2.0	ND	2.0	ND	2.0	ND	2.0	ND	2.0
Chloroethane		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Trichlorofluoromet	hane	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Acetone		ND	2.0	6.2	2.0	3.4	2.0	7.5	2.0	3.7	2.0
1,1-Dichloroethene	}	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Carbon Disulfide		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Methylene Chloride	•	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Methyl-tertiary-bu	itylether	ND	1.0	ND	1.0	14	1.0	ND	1.0	ND	1.0
trans-1,2-Dichloro	ethene	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,1-Dichloroethane		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
2-Butanone		ND	4.0	ND	4.0	ND	4.0	ND	4.0	ND	4.0
2,2-Dichloropropan	e	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
cis-1,2-Dichloroet	hene	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Chloroform		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,1-Dichloropropen	e	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,2-Dichloroethane		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,1,1-Trichloroeth	ane	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Carbon Tetrachlori		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Benzene		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Trichloroethene		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,2-Dichtoropropan	e	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Dibromomethane		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Bromodichlorometha	ne	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
cis-1,3-Dichloropr	opene	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
trans-1,3-Dichloro	propene	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,1,2-Trichloroeth	ane	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,3-Dichloropropan	e	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Dibromochtorometha	ne	ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,2-Dibromoethane		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Bromoform		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
4-Methyl-2-Pentano	ne	ND	2.0	ND	2.0	ND	2.0	ND	2.0	ND	2.0
-											

Table 1.1 (Cont) Results of the VOA Analysis for the Water Samples WA # 0-098 Pier Drum

	SAMPLE # LOCATION COLLECTED ANALYZED INJECTED FILE DIL. FACT.	:	01/18/95 00:24 ^V2472	K	00636 MW-S DUP 01/11/95 01/18/95 01:04 ^V2473 1	i	00641 MM-N 01/11/95 01/18/95 01:45 ^V2474 1		00635 MW-S 01/11/95 01/18/95 02:25 ^V2475 1		00634 MW-W 01/11/95 01/18/95 03:46 ^V2477	
	UNIT	:	ug/L		ug/L		ug/L		ug/L		ug/L	
COMPOUND			CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL
Toluene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
2-Hexanone			ND	2.0	ND	2.0	ND	2.0	ND	2.0	ND	2.0
Tetrachloroethene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Chlorobenzene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,1,1,2-Tetrachlord	pethane		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Ethylbenzene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
p & m-Xylene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
o-Xylene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Styrene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Isopropylbenzene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,1,2,2-Tetrachlord	oethane		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,2,3-Trichloroprop	oane		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Bromobenzene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
n-Propylbenzene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
2-Chlorotoluene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
4-Chiorotoluene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,3,5-Trimethylbenz	ene		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
tert-Butylbenzene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,2,4-Trimethylbenz	ene		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
sec-Butylbenzene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,3-Dichlorobenzene	•		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
p-Isopropyltoluene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,4-Dichlorobenzene	•		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,2-Dichlorobenzene	•		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
n-Butylbenzene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1,2-Dibromo-3-Chlor	opropane		ND	1.0	MD	1.0	ND	1.0	ND	1.0	ND	1.0
1,2,4-Trichlorobenz	ene		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Naphthalene			ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
Hexachlorobutadiene	•		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0
1.2.3-Trichlorobenz	ene		ND	1.0	ND	1.0	ND	1.0	ND	1.0	ND	1.0

Table 1.1 (Cont) Results of the VOA Analysis for the Water Samples WA # 0-098 Pier Drum

SAMPLE #	:	LAB BLANK		00637
LOCATION	:			MW-E
COLLECTED	:			01/11/95
ANALYZED :	:	01/18/95		01/18/95
INJECTED :	:	00:24		04:27
FILE #	:	^V2472		^v2478
DIL. FACT.:	:	1		1
UNIT	:	ug/L		ug/L
		CONC.	MDL	CONC.

COMPOUND	CONC.	MDL	CONC.	MDL
Dichlorodifluoromethane	ND	1.0	ND	1.0
Chloromethane	ND	1.0	ND	1.0
Vinyl Chloride	ND	1.0	ND	1.0
Bromomethane	ND	2.0	ND	2.0
Chloroethane	ND	1.0	ND	1.0
Trichlorofluoromethane	ND	1.0	ND	1.0
Acetone	ND	2.0	8.0	2.0
1,1-Dichloroethene	ND	1.0	ND	1.0
Carbon Disulfide	ND	1.0	ND	1.0
Methylene Chloride	ND	1.0	ND	1.0
Methyl-tertiary-butylether	ND	1.0	ND	1.0
trans-1,2-Dichloroethene	ND	1.0	ND	1.0
1,1-Dichloroethane	ND	1.0	ND	1.0
2-Butanone	ND	4.0	0.8 J	4.0
2,2-Dichloropropane	ND	1.0	ND	1.0
cis-1,2-Dichloroethene	ND	1.0	ND	1.0
Chloroform	ND	1.0	ND	1.0
1,1-Dichloropropene	ND	1.0	ND	1.0
1,2-Dichloroethane	ND	1.0	ND	1.0
1,1,1-Trichloroethane	ND	1.0	ND	1.0
Carbon Tetrachloride	ND	1.0	ND	1.0
8enzene	ND	1.0	ND	1.0
Trichloroethene	ND	1.0	ND	1.0
1,2-Dichloropropane	ND	1.0	ND	1.0
Dibromomethane	ND	1.0	ND	1.0
Bromodichloromethane	ND	1.0	ND	1.0
cis-1,3-Dichloropropene	ND	1.0	ND	1.0
trans-1,3-Dichloropropene	ND	1.0	ND	1.0
1,1,2-Trichloroethane	ND	1.0	ND	1.0
1,3-Dichloropropane	ND	1.0	ND	1.0
Dibromochloromethane	ND	1.0	ND	1.0
1,2-Dibromoethane	ND	1.0	ND	1.0
Bromoform	ND	1.0	ND	1.0
4-Methyl-2-Pentanone	ND	2.0	ND	2.0

Table 1.1 (Cont) Results of the VOA Analysis for the Water Samples WA # 0-098 Pier Drum

SAMPLE # :	LAB BLANK	00637
LOCATION :		MW-E
COLLECTED :		01/11/95
ANALYZED :	01/18/95	01/18/95
INJECTED :	00:24	04:27
FILE :	^V2472	^v2478
DIL. FACT .:	1	1
UNIT :	ug/L	ug/L

COMPOUND	CONC.	MDL	CONC.	MDL
Toluene	ND	1.0	2.2	1.0
2-Hexanone	ND	2.0	ND	2.0
Tetrachloroethene	ND	1.0	ND	1.0
Chlorobenzene	ND	1.0	ND	1.0
1,1,1,2-Tetrachloroethane	ND	1.0	ND	1.0
Ethylbenzene	ND	1.0	ND	1.0
p & m-Xylene	ND	1.0	ND	1.0
o-Xylene	ND	1.0	ND	1.0
Styrene	ND	1.0	ND	1.0
1sopropylbenzene	ND	1.0	ND	1.0
1,1,2,2-Tetrachloroethane	ND	1.0	ND	1.0
1,2,3-Trichloropropane	ND	1.0	ND	1.0
Bromobenzene	ND	1.0	ND	1.0
n-Propylbenzene	ND	1.0	ND	1.0
2-Chlorotoluene	ND	1.0	ND	1.0
4-Chiorotoluene	ND	1.0	ND	1.0
1,3,5-Trimethylbenzene	ND	1.0	ND	1.0
tert-Butylbenzene	ND	1.0	ND	1.0
1,2,4-Trimethylbenzene	ND	1.0	ND	1.0
sec-Butylbenzene	ND	1.0	ND	1.0
1,3-Dichlorobenzene	ND	1.0	ND	1.0
p-Isopropyltoluene	ND	1.0	ND	1.0
1,4-Dichlorobenzene	ND	1.0	ND	1.0
1,2-Dichlorobenzene	ND	1.0	ND	1.0
n-Butylbenzene	ND	1.0	ND	1.0
1,2-Dibromo-3-Chloropropane	ND	1.0	ND	1.0
1,2,4-Trichlorobenzene	ND	1.0	ND	1.0
Naphthalene	ND	1.0	ND	1.0
Hexachlorobutadiene	ND	1.0	ND	1.0
1,2,3-Trichlorobenzene	ND	1.0	ND	1.0

Table 1.1 (Cont) Results of the VOA Analysis for the Water Samples WA # 0-098 Pier Drum

SAMPLE #	:	LAB BLANK	00639
LOCATION	:		W PIT
COLLECTED	:		01/11/95
ANALYZED	:	01/18/95	01/18/95
INJECTED	:	09:48	11:09
FILE #	:	^V2481	^V2483
DIL. FACT.	:	1	1
UNIT	:	ua/L	ua/L

COMPOUND	CONC.	MDL	CONC.	MDL
Dichlorodifluoromethane	ND	1.0	ND	1.0
Chloromethane	ND	1.0	ND	1.0
Vinyl Chloride	ND	1.0	ND	1.0
Bromomethane	ND	2.0	ND	2.0
Chloroethane	ND	1.0	ND	1.0
Trichlorofluoromethane	ND	1.0	ND	1.0
Acetone	ND	2.0	7.3	2.0
1,1-Dichloroethene	ND	1.0	ND	1.0
Carbon Disulfide	ND	1.0	ND	1.0
Methylene Chloride	ND	1.0	ND	1.0
Methyl-tertiary-butylether	ND	1.0	ND	1.0
trans-1,2-Dichloroethene	ND	1.0	ND	1.0
1,1-Dichloroethane	ND	1.0	ND	1.0
2-Butanone	ND	4.0	ND	4.0
2,2-Dichloropropane	ND	1.0	ND	1.0
cis-1,2-Dichloroethene	ND	1.0	ND	1.0
Chloroform	ND	1.0	ND	1.0
1,1-Dichloropropene	ND	1.0	ND	1.0
1,2-Dichloroethane	ND	1.0	ND	1.0
1,1,1-Trichloroethane	ND	1.0	ND	1.0
Carbon Tetrachioride	ND	1.0	ND	1.0
Benzene	ND	1.0	ND	1.0
Trichloroethene	ND	1.0	ND	1.0
1,2-Dichloropropane	ND	1.0	ND	1.0
Dibromomethane	ND	1.0	ND	1.0
Bromodichioromethane	ND	1.0	ND	1.0
cis-1,3-Dichloropropene	ND	1.0	ND	1.0
trans-1,3-Dichloropropene	ND	1.0	ND	1.0
1,1,2-Trichloroethane	ND	1.0	ND	1.0
1,3-Dichloropropane	ND	1.0	ND	1.0
Dibromochloromethane	ND	1.0	ND	1.0
1,2-Dibromoethane	ND	1.0	ND	1.0
Bromoform	ND	1.0	ND	1.0
4-Methyl-2-Pentanone	ND	2.0	ND	2.0

Table 1.1 (Cont) Results of the VOA Analysis for the Water Samples WA # 0-098 Pier Drum

SAMPLE # :	LAB BLANK	00639
LOCATION :		W PIT
COLLECTED :		01/11/95
ANALYZED :	01/18/95	01/18/95
INJECTED :	09:48	11:09
FILE :	^V2481	^V2483
DIL. FACT .:	1	1
UNIT :	ua/L	ua/L

•			-g, -	
COMPOUND	CONC.	MDL	CONC.	MDL
Toluene	ND	1.0	ND	1.0
2-Hexanone	ND	2.0	ND	2.0
Tetrachloroethene	ND	1.0	ND	1.0
Chlorobenzene	ND	1.0	ND	1.0
1,1,1,2-Tetrachloroethane	ND	1.0	ND	1.0
Ethylbenzene	ND	1.0	ND	1.0
p & m-Xylene	ND	1.0	ND	1.0
o-Xylene	ND	1.0	ND	1.0
Styrene	ND	1.0	ND	1.0
lsopropylbenzene	ND	1.0	ND	1.0
1,1,2,2-Tetrachloroethane	ND	1.0	ND	1.0
1,2,3-Trichloropropane	ND	1.0	ND	1.0
Bromobenzene	ND	1.0	ND	1.0
n-Propylbenzene	ND	1.0	ND	1.0
2-Chlorotoluene	ND	1.0	ND	1.0
4-Chlorotoluene	ND	1.0	ND	1.0
1,3,5-Trimethylbenzene	ND	1.0	ND	1.0
tert-Butylbenzene	ND	1.0	ND	1.0
1,2,4-Trimethylbenzene	ND	1.0	ND	1.0
sec-Butylbenzene	ND	1.0	ND	1.0
1,3-Dichlorobenzene	ND	1.0	ND	1.0
p-Isopropyltoluene	ND	1.0	ND	1.0
1,4-Dichlorobenzene	ND	1.0	ND	1.0
1,2-Dichlorobenzene	ND	1.0	ND	1.0
n-Butylbenzene	ND	1.0	ND	1.0
1,2-Dibromo-3-Chloropropane	ND	1.0	ND	1.0
1,2,4-Trichlorobenzene	ND	1.0	ND	1.0
Naphthalene	ND	1.0	ND	1.0
Hexachlorobutadiene	ND	1.0	ND	1.0
1,2,3-Trichlorobenzene	ND	1.0	ND	1.0

Table 1.2 Results of the VOA in Water TICs WA # 0-098 Pier Drum

Sample ID	Location	Compound	Concentration* (ug/l)	Retention Time (minutes)
Lab Blank 1/13/95		Alkane	7	23.66
00638	N Pit	None Found		
00645	Trip Blank	None Found		
Lab Blank 1/14/95		None Found		
00640	Field blank	None Found		
Lab Blank 1/18/95		None Found		
00636	MW-Dup	None Found		
00641	MW-N	None Found		
00635	MW-S	None Found		
00634	MW-W	None Found		
00637	MW-E	None Found		
Lab Blank 1/18/95		None Found		
00639	West Pit (WPit)	None Found		

^{*} denotes that the concentration is estimated - the response factor was assumed to be 1.

Table 1.3 Results of the VOA Analysis for the Soil Samples $$\rm WA~\#~0\text{-}098~Pier~Drum$$

17 S

	SAMPLE # : LOCATION : COLLECTED : ANALYZED : INJECTED : FILE # : DIL. FACT.: % SOLID : UNIT :	01/14/9: 13:30 ^V2442 1 100 ug/Kg		00643 SED N 01/11/9: 01/14/9: 19:35 ^V2451 1 81 ug/Kg		00644 SED E 01/11/9 01/14/9 20:16 ^V2452 1 79 ug/Kg	
COMPOUND		CONC.	MDL	CONC.	MDL	CONC.	MDL
Dichlorodifluorome	thane	ND	1.0	ND	1.2	ND	1.3
Chloromethane		ND	1.0	ND	1.2	ND	1.3
Vinyl Chloride		ND	1.0	ND	1.2	ND	1.3
Bromomethane		ND	2.0	ND	2.5	ND	2.5
Chloroethane		ND	1.0	ND	1.2	ND	1.3
Trichlorofluoromet	han e	ND	1.0	ND	1.2	ND	1.3
Acetone		ND	2.0	ND	2.5	ND	2.5
1,1-Dichloroethene		ND	1.0	ND	1.2	ND	1.3
Carbon Disulfide		ND	1.0	ND	1.2	ND	1.3
Methylene Chloride	!	ND	1.0	ND	1.2	ND	1.3
Methyl-tertiary-bu		ND	1.0	ND	1.2	ND	1.3
trans-1,2-Dichloro	ethene	ND	1.0	ND	1.2	ND	1.3
1,1-Dichloroethane	•	ND	1.0	ND	1.2	ND	1.3
2-Butanone		ND	4.0	ND	4.9	ND	5.1
2,2-Dichloropropan	e	ND	1.0	ND	1.2	ND	1.3
cis-1,2-Dichloroet	hene	ND	1.0	ND	1.2	ND	1.3
Chloroform		ND	1.0	ND	1.2	ND	1.3
1,1-Dichloropropen	e	ND	1.0	ND	1.2	ND	1.3
1,2-Dichloroethane		ND	1.0	ND	1.2	ND	1.3
1,1,1-Trichloroeth		ND	1.0	ND	1.2	ND	1.3
Carbon Tetrachlori	de	ND	1.0	ND	1.2	ND	1.3
Benzene		ND	1.0	ND	1.2	ND	1.3
Trichloroethene		ND	1.0	ND	1.2	ND	1.3
1,2-Dichloropropan	e	ND	1.0	ND	1.2	ND	1.3
Dibromomethane		ND	1.0	ND	1.2	ND	1.3
Bromodichlorometha	ne	ND	1.0	ND	1.2	ND	1.3
cis-1,3-Dichloropr		ND	1.0	ND	1.2	ND	1.3
trans-1,3-Dichloro		ND	1.0	ND	1.2	ND	1.3
1,1,2-Trichloroeth		ND	1.0	ND	1.2	ND	1.3
1,3-Dichloropropan		ND	1.0	ND	1.2	ND	1.3
Dibromochlorometha	ne	ND	1.0	ND	1.2	ND	1.3
1,2-Dibromoethane		ND	1.0	ND	1.2	ND	1.3
Bromoform		ND	1.0	ND	1.2	ND	1.3
4-Methyl-2-Pentano	ne	ND	2.0	ND	2.5	ND	2.5

Table 1.3 (Cont) Results of the VDA Analysis for the Soil Samples
WA # 0-098 Pier Drum

LOCA COLL ANAL INJE FILE DIL	. FACT.: OLID :	01/14/95 13:30 ~V2442 1 100 ug/Kg	•	00643 SED N 01/11/95 01/14/95 19:35 ^V2451 1 81 ug/Kg	S 0 0 2 2	0644 ED E 1/11/95 1/14/95 0:16 V2452 1 79 ug/Kg	MDL
COMPOUND		CONC.	MDL	CONC.	MDL		1.3
Toluene 2-Hexanone Tetrachloroethene Chlorobenzene 1,1,1,2-Tetrachloroetl Ethylbenzene p & m-Xylene o-Xylene Styrene Isopropylbenzene 1,2,2-Tetrachloroetl 1,2,3-Trichloropropar Bromobenzene n-Propylbenzene 2-Chlorotoluene 4-Chlorotoluene 1,3,5-Trimethylbenzene tert-Butylbenzene 1,2,4-Trimethylbenzene 1,3-Dichlorobenzene p-Isopropyltoluene 1,4-Dichlorobenzene 1,2-Dichlorobenzene 1,2-Dichlorobenzene 1,2-Dichlorobenzene 1,2-Trichlorobenzene 1,2,4-Trichlorobenzene 1,2,4-Trichlorobenzene 1,2,3-Trichlorobenzene	thane ne ene ene eopropane zene	ND N	1.0 2.0 1.0 1.0 1.0 1.0 1.0 1.0 1.0 1	ND ND ND ND ND ND ND ND	1.2 2.5 1.2 1.2 1.2 1.2 1.2 1.2 1.2 1.2 1.2 1.2	ND ND ND ND	1.3 1.3 1.3 1.3 1.3 1.3 1.3 1.3 1.3 1.3

3343

Table 1.3 (Cont) Results of the VOA Analysis for the Soil Samples WA # 0-098 Pier Drum

	SAMPLE # :	LAB BLA	NK	00642	
	LOCATION :			SED-S	
	COLLECTED :			01/11/9	5
	ANALYZED :	01/17/9	5	01/17/9	5
	INJECTED :	15:41		16:35	
	FILE # :	^V2460		^V2461	
	DIL. FACT .:	1		1	
	% SOLID :	100		79	
	UNIT :	ug/Kg		ug/Kg	
				J. J	
COMPOUND		CONC.	MDL	CONC.	MDL
Dichlorodifluoromet	hane	ND	1.0	ND	1.3
Chloromethane		ND	1.0	ND	1.3
Vinyl Chloride		ND	1.0	ND	1.3
Bromomethane		ND	2.0	ND	2.5
Chloroethane		ND	1.0	ND	1.3
Trichlorofluorometh	ane	ND	1.0	ND	1.3
Acetone	J. 1.C	ND	2.0	207	2.5
1.1-Dichloroethene		ND	1.0	ND	1.3
Carbon Disulfide		ND	1.0	ND	1.3
Methylene Chloride		ND	1.0	ND	1.3
Methyl-tertiary-but	dathar	ND	1.0	ND	1.3
trans-1,2-Dichloroe		ND	1.0	ND	1.3
· ·	Literie	ND	1.0	ND	1.3
1,1-Dichloroethane		ND	4.0	ND	5.1
2-Butanone			1.0		1.3
2,2-Dichloropropane		ND		ND	
cis-1,2-Dichloroeth	ene	ND	1.0	ND	1.3
Chloroform		ND	1.0	ND	1.3
1,1-Dichloropropene		ND	1.0	ND	1.3
1,2-Dichloroethane		ND	1.0	ND	1.3
1,1,1-Trichloroetha		ND	1.0	ND	1.3
Carbon Tetrachloride	2	ND	1.0	ND	1.3
Benzene		ND	1.0	ND	1.3
Trichloroethene		ND	1.0	ND	1.3
1,2-Dichloropropane		ND	1.0	ND	1.3
Dibromomethane		ND	1.0	ND	1.3
Bromodichtoromethan		ND	1.0	ND	1.3
cis-1,3-Dichloroprop		ND	1.0	ND	1.3
trans-1,3-Dichlorop		ND	1.0	ND	1.3
1,1,2-Trichloroetha	ne	ND	1.0	ND	1.3
1,3-Dichtoropropane		ND	1.0	ND	1.3
Dibromochloromethan	2	ND	1.0	ND	1.3
1,2-Dibromoethane		ND	1.0	ND	1.3
Bromoform		ND	1.0	ND	1.3
4-Methyl - 2-Pentanon	9	MD	2 0	NO	2.5

ND

1.0

ND

1.3

4-Methyl-2-Pentanone

Table 1.3 (Cont) Results of the VOA Analysis for the Soil Samples WA # 0-098 Pier Drum

SAMPLE #	:	LAB BLANK	00642
LOCATION	:		SED-S
COLLECTED	:		01/11/95
ANALYZED	:	01/17/95	01/17/95
INJECTED	:	15:41	16:35
FILE	:	^V2460	^V2461
DIL. FACT.	. :	1	1
% SOLID	:	100	79
UNIT	:	ug/Kg	ug/Kg

COMPOUND	CONC.	MDL	CONC.	MDL
Toluene	ND	1.0	12	1.3
2-Hexanone	ND	2.0	ND	2.5
Tetrachloroethene	ND	1.0	ND	1.3
Chlorobenzene	ND	1.0	ND	1.3
1,1,1,2-Tetrachloroethane	ND	1.0	ND	1.3
Ethylbenzene	ND	1.0	670	1.3
p & m-Xylene	ND	1.0	ND	1.3
o-Xylene	ND	1.0	2.4	1.3
Styrene	ND	1.0	2.8	1.3
Isopropylbenzene	ND	1.0	2.1	1.3
1,1,2,2-Tetrachloroethane	ND	1.0	ND	1.3
1,2,3-Trichloropropane	ND	1.0	ND	1.3
Bromobenzene	ND	1.0	ND	1.3
n-Propylbenzene	ND	1.0	1.6	1.3
2-Chlorotoluene	ND	1.0	ND	1.3
4-Chlorotoluene	ND	1.0	ND	1.3
1,3,5-Trimethylbenzene	ND	1.0	ND	1.3
tert-Butylbenzene	ND	1.0	ND	1.3
1,2,4-Trimethylbenzene	ND	1.0	3.3	1.3
sec-Butylbenzene	ND	1.0	ND	1.3
1,3-Dichlorobenzene	ND	1.0	ND	1.3
p-Isopropyltoluene	ND	1.0	1.2 J	1.3
1,4-Dichlorobenzene	ND	1.0	ND	1.3
1,2-Dichlorobenzene	ND	1.0	ND	1.3
n-Butylbenzene	ND	1.0	ND	1.3
1,2-Dibromo-3-Chloropropane	ND	1.0	ND	1.3
1,2,4-Trichlorobenzene	ND	1.0	ND	1.3
Naphthalene	ND	1.0	ND	1.3
Hexach Lorobutadiene	ND	1.0	ND	1.3
1,2,3-Trichlorobenzene	ND	1.0	ND	1.3

2 2 35.0

Table 1.4 Results of the VOA in Soil TICs WA # 0-098 Pier Drum

Sample ID	Location	Compound	Concentration*	Retention Time	
			(ug/l)	(minutes)	
Lab Blank 1/14/95		None Found			
00643	Sed N	None Found			
00644	Sed E	None Found			
Lab Blank 1/17/95		None Found	1		
00642	Sed S	Alkyl Benzene	8	20.67	

^{*} denotes that the concentration is estimated - the response factor was assumed to be 1.

Table 1.5 Results of the TCLP VOA Analysis WA # 0-098 Pier Drum

Sample # Location Collected		ratory ank		0644 d E	B 00 Sec		B 00 Sec	
Analysed	_		_		_		_	
Dil. Factor	Conc	MDL	Conc	MDL	Conc	MDL	Conc	MDL
Compound	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l
Chloromethane	ND	10	ND	100	ND	100	ND	100
Bromomethane	ND	10	ND	100	ND	100	ND	100
Vinyl chloride	ND	10	ND	100	ND	100	ND	100
Chloroethane	ND	10	ND	100	ND	100	ND	100
Methylene chloride	2.3	J 10	ND	100	ND	100	ND	100
Acetone	ND	50	ND	50 0	ND	500	ND	500
Carbon Disulfide	ND	10	ND	100	ND	100	ND	100
1,1-Dichloroethene	ND	10	ND	100	ND	100	ND	100
1,1-Dichloroethane	ND	10	ND	100	ND	100	ND	100
cis/trans-1,2-Dichloroethene	ND	10	ND	100	ND	100	ND	100
Chloroform	ND	10	ND	100	ND	100	ND	100
1,2-Dichloroethane	ND	10	ND	100	ND	100	ND	100
2-Butanone	ND	50	ND	500	ND	500	ND	500
1,1,1,~Trichloroethane	ND	10	ND	100	ND	100	ND	100
Carbon tetrachloride	ND	10	ND	100	ND	100	ND	100
Vinyl Acetate	ND	50	ND	500	ND	500	ND	500
Bromodichloromethane	ND	10	ND	100	ND	100	ND	100
1,2-Dichloropropane	ND	10	ND	100	ND	100	ND	100
Trichloroethene	ND	10	ND	100	ND	100	ND	100
Benzene	ND	10	ND	100	ND	100	ND	100
cis-1,3-Dichloropropene	ND	10	ND	100	ND	100	ND	100
Dibromochloromethane	ND	10	ND	100	ND	100	ND	100
trans-1,3-Dichloropropene	ND	10	ND.	100	ND	100	ND	100
1,1,2-Trichloroethane	ND	10	ND	100	ND	100	ND	100
Bromoform	ND	10	ND	100	ND	100	ND	100
4-Methyl-2-Pentanone	ND	50	ND	500	ND	500	ND	500
2-Hexanone	ND	50	ND	500	ND	500	ND	500
1.1.2.2 - Tetrachloroethane	ND	10	ND	100	ND	100	ND	100
Tetrachloroethene	ND	10	ND	100	ND	100	ND	100
Toluene	1.3	J 10	ND	100	ND	100	ND	100
Chlorobenzene	ND	10	ND	100	ND	100	ND	100
Ethyl benzene	ND	10	ND	100	ND	100	ND	100
Styrene	ND	10	ND	100	ND	100	ND	100
m-Xylene	ND	10	ND	100	ND	100	ND	100
o/p-Xylene	ND	20	ND	200	ND	200	ND	200
or programme	.,						,	

Table 1.6 Results of the VOA Analysis for the Drum Liquid Samples WA # 0-098 Pier Drum

SAMPLE #	:	LAB BLANK	00914
LOCATION	:		REAC #4
COLLECTED	:		01/10/95
ANALYZED	:	01/30/95	01/30/95
INJECTED		17:11	17:52
FILE #	:	^V2619	^V2620
DIL. FACT.	:	1	1
UNIT :	:	ug/L	ug/L

UNII	•	ug/L		ug/L	
COMPOUND		CONC.	MDL	CONC.	MDL
Dichlorodifluoromethane		ND	1.0	ND	1.0
Chloromethane		ND	1.0	ND	1.0
Vinyl Chloride		ND	1.0	ND	1.0
Bromomethane		ND	2.0	ND	2.0
Chloroethane		ND	1.0	ND	1.0
Trichlorofluoromethane		ND	1.0	ND	1.0
Acetone		ND	2.0	1400	2.0
1,1-Dichloroethene		ND	1.0	ND	1.0
Carbon Disulfide		ND	1.0	ND	1.0
Methylene Chloride		ND	1.0	ND	1.0
Methyl-tertiary-butylether		ND	1.0	ND	1.0
trans-1,2-Dichloroethene		ND	1.0	ND	1.0
1,1-Dichloroethane		ND	1.0	ND	1.0
2-Butanone		ND	4.0	89	4.0
2,2-Dichloropropane		ND	1.0	ND	1.0
cis-1,2-Dichloroethene		ND	1.0	ND	1.0
Chloroform		ND	1.0	ND	1.0
1,1-Dichloropropene		ND	1.0	ND	1.0
1,2-Dichloroethane		ND	1.0	ND	1.0
1,1,1-Trichloroethane		ND	1.0	ND	1.0
Carbon Tetrachloride		ND	1.0	ND	1.0
Benzene		ND	1.0	ND	1.0
Trichloroethene		ND	1.0	ND	1.0
1,2-Dichloropropane		ND	1.0	ND	1.0
Dibromomethane		ND	1.0	ND	1.0
Bromodichloromethane		ND	1.0	ND	1.0
cis-1,3-Dichloropropene		ND	1.0	ND	1.0
trans-1,3-Dichloropropene		ND	1.0	ND	1.0
1,1,2-Trichloroethane		ND	1.0	ND	1.0
1,3-Dichloropropane		ND	1.0	ND	1.0
Dibromochloromethane		ND	1.0	ND	1.0
1,2-Dibromoethane		ND	1.0	ND	1.0
Bromoform		ND	1.0	ND	1.0
4-Methyl-2-Pentanone		ND	2.0	ND	2.0

2 2 3313

Table 1.6 (Cont) Results of the VOA Analysis for the Drum Liquid Samples WA # 0-098 Pier Drum

SAMPLE # :	LAB BLANK	00914
LOCATION :		REAC #4
COLLECTED :		01/10/95
ANALYZED :	01/30/95	01/30/95
INJECTED :	17:11	17:52
fILE :	^V2619	^v2620
DIL. FACT .:	1	1
UNIT :	ug/L	ug/L

ONII	. Ug/L		ug/L	
COMPOUND	CONC.	MDL	CONC.	MDL
Toluene	ND	1.0	9.6	1.0
2-Hexanone	ND	2.0	ND	2.0
Tetrachloroethene	ND	1.0	ND	1.0
Chlorobenzene	ND	1.0	ND	1.0
1,1,1,2-Tetrachloroethane	ND	1.0	ND	1.0
Ethylbenzene	ND	1.0	37	1.0
p & m-Xylene	ND	1.0	ND	1.0
o-Xylene	ND	1.0	0.8 J	1.0
Styrene	ND	1.0	26	1.0
Isopropylbenzene	ND	1.0	ND	1.0
1,1,2,2-Tetrachloroethane	ND	1.0	ND	1.0
1,2,3-Trichloropropane	ND	1.0	ND	1.0
Bromobenzene	ND	1.0	ND	1.0
n-Propylbenzene	ND	1.0	ND	1.0
2-Chlorotoluene	ND	1.0	ND	1.0
4-Chlorotoluene	ND	1.0	ND	1.0
1,3,5-Trimethylbenzene	ND	1.0	ND	1.0
tert-Butylbenzene	ND	1.0	ND	1.0
1,2,4-Trimethylbenzene	ND	1.0	45	1.0
sec-Butylbenzene	ND	1.0	ND	1.0
1,3-Dichlorobenzene	ND	1.0	ND	1.0
p-Isopropyltoluene	ND	1.0	ND	1.0
1,4-Dichlorobenzene	ND	1.0	ND	1.0
1,2-Dichlorobenzene	ND	1.0	ND	1.0
n-Butylbenzene	ND	1.0	ND	1.0
1,2-Dibromo-3-Chloropropane	ND	1.0	ND	1.0
1,2,4-Trichlorobenzene	ND	1.0	ND	1.0
Naphthalene	ND	1.0	7.2	1.0
Hexach Lorobutadiene	ND	1.0	ND	1.0
1,2,3-Trichtorobenzene	ND	1.0	ND	1.0

2 () 3314

Table 1.6 (Cont) Results of the VOA Analysis for the Drum Liquid Samples WA # 0-098 Pier Drum

	SAMPLE # : LOCATION : COLLECTED : ANALYZED : INJECTED : FILE # : DIL. FACT.: UNIT :	01/30/95 17:11 ^V2619 1		00912 REAC #9 01/10/9 01/33/9 18:33 ^V2621 5 mg/L	5	00913 REAC #2 01/10/9 01/30/9 19:14 ^V2622 5 mg/L	95
COMPOUND		CONC.	MDL	CONC.	MD	CONC.	MDL
Dichlorodifluorome Chloromethane Vinyl Chloride Bromomethane Chloroethane Trichlorofluorometh Acetone 1,1-Dichloroethene Carbon Disulfide Methylene Chloride T,2-Dichloroethane 2,2-Dichloropropane cis-1,2-Dichloroeth Chloroform 1,1-Dichloropropane 1,2-Dichloroethane	nane tylether ethene e nene	ND N	1.0 1.0 1.0 1.0 1.0 4.0 1.0 1.0 1.0	ND ND ND ND ND 130000 ND ND ND ND ND ND ND ND ND ND ND	5.0 5.0 5.0 10.0 5.0 5.0 5.0 5.0 5.0 5.0 5.0 5.0 5.0	ND N	5.0 5.0 5.0 10.0 5.0 10.0 5.0 5.0 5.0 5.0 5.0 5.0 5.0 5.0
1,1,1-Trichloroetha Carbon Tetrachloric		ND ND	1.0	ND ND	5.0 5.0	ND ND	5.0 5.0
Benzene Trichloroethene 1,2-Dichloropropane Dibromomethane		ND ND ND ND	1.0 1.0 1.0	11 ND ND ND	5.0 5.0 5.0 5.0	12 ND ND ND	5.0 5.0 5.0 5.0
Bromodichloromethan cis-1,3-Dichloropro trans-1,3-Dichlorop 1,1,2-Trichloroetha	opene oropene ane	ND ND ND ND	1.0 1.0 1.0	ND ND ND ND	5.0 5.0 5.0	ND ND ND ND	5.0 5.0 5.0
1,3-Dichloropropane Dibromochloromethan 1,2-Dibromoethane Bromoform 4-Methyl-2-Pentanor	ne	ND ND ND ND ND	1.0 1.0 1.0 1.0 2.0	ND ND ND ND ND	5.0 5.0 5.0 5.0 10.0	ND ND ND ND ND	5.0 5.0 5.0 5.0

2 2 23 5

Table 1.6 (Cont) Results of the VOA Analysis for the Drum Liquid Samples $$\rm WA~\#~0\text{-}098~Pier~Drum$$

SAMPLE # : LOCATION : COLLECTED : ANALYZED : INJECTED : FILE : DIL. FACT.: UNIT :	01/30/9 17:11 ^V2619 1		00912 REAC #9 01/10/9 01/30/9 18:33 ^V2621 5 mg/L	5	00913 REAC #2 01/10/9 01/30/9 19:14 ^V2622 5 mg/L	5
COMPOUND	CONC.	MDL	CONC.	MDL	CONC.	MDL
Toluene	ND	1.0	5.4	5.0	4.1 J	5.0
2-Hexanone	ND	2.0	ND	10.0	ND	10.0
Tetrachloroethene	ND	1.0	ND	5.0	ND	5.0
Chlorobenzene	ND	1.0	ND	5.0	ND	5.0
1,1,1,2-Tetrachioroethane	ND	1.0	ND	5.0	ND	5.0
Ethylbenzene	ND	1.0	21	5.0	88	5.0
p & m-Xylene	ND	1.0	16	5.0	51	5.0
o-Xylene	ND	1.0	8.4	5.0	ND	5.0
Styrene	ND	1.0	1200		16000	5.0
Isopropylbenzene	ND	1.0	6.5	5.0	21	5.0
1,1,2,2-Tetrachloroethane	ND	1.0	ND	5.0	ND	5.0
1,2,3-Trichloropropane	ND	1.0	ND	5.0	ND	5.0
Bromobenzene	ND	1.0	ND	5.0	ND	5.0
n-Propylbenzene	ND	1.0	7.6	5.0	17	5.0
2-Chlorotoluene	ND	1.0	ND	5.0	ND	5.0
4-Chlorotoluene	ND	1.0	ND	5.0	ND	5.0
1,3,5-Trimethylbenzene	ND	1.0	ND	5.0	ND	5.0
tert-Butylbenzene	ND	1.0	ND	5.0	ND	5.0
1,2,4-Trimethylbenzene	ND	1.0	17	5.0	23	5.0
sec-Butylbenzene	ND ND	1.0	ND	5.0	ND	5.0
1,3-Dichlorobenzene	ND	1.0	ND	5.0 5.0	ND	5.0
p-Isopropyltoluene	ND ND	1.0 1.0	ND ND	5.0	ND ND	5.0 5.0
1,4-Dichlorobenzene	ND	1.0	ND	5.0	ND	5.0
1,2-Dichlorobenzene	ND	1.0	ND	5.0	ND	5.0
n-Butylbenzene	ND	1.0	ND	5.0	ND	5.0
1,2-Dibromo-3-Chloropropane	ND	1.0	ND	5.0	ND	5.0
1,2,4-Trichlorobenzene Naphthalene	ND	1.0	3.4 J	5.0	ND	5.0
Hexachlorobutadiene	ND	1.0	ND	5.0	ND	5.0
1,2,3-Trichlorobenzene	ND	1.0	ND	5.0	ND	5.0
1,2,3 IT followbenzene	NU	1.0	NU	3.0	NU	٥.٠

2 2 9316

Table 1.7 Results of the VOA in Drum Liquid TICs WA # 0-098 Pier Drum

Sample ID	Location	Compound	Concentration* (ug/l)	Retention Time (minutes)
Lab Blank 1/30/95		None Found		
00914	REAC #4	Ketone	8	18.39
00912	REAC #9	Unknown Alkyl Benzene	160 30	5.94 28.68
00913	REAC #2	Unknown Alkyl Benzene Unknown	170 75 50	5.91 20.68 21.42
Lab Blank 1/31/95		None Found		

^{*} denotes that the concentration is estimated - the response factor was assumed to be 1.

Table 1.8 Results of the TCLP VOA Analysis for the Drum Solid Samples WA # 0-098 Pier Drum

Sample # Location Collected Analysed	Lab Blank		00911 Composit 1, 8, 9, 11		00915 REAC #5		00916 Composit 3, 7, 10, 13, 14	1
Dil. Factor	Conc	MDL	Conc	MDL	Conc	MDL	Conc	MDL
Compound	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	ug/i	ug/l
Chloromethane	ND	10	ND	5000	ND	5000	ND	2500
Bromomethane	ND	10	ND	5000	ND	5000	ND	2500
Vinyl chloride	ND	10	ND	5000	ND	5000	ND	2500
Chloroethane	ND	10	ND	5000	ND	5000	ND	2500
Methylene chloride	2.3		1300 JB	5000	1100 JB		ND	2500
Acetone	ND	- 10	1300000	5000	11000 J	5000	ND	2500
Carbon Disulfide	ND	10	ND	5000	ND	5000	ND	2500
1,1-Dichloroethene	ND	10	ND	5000	ND	5000	ND	2500
1,1-Dichloroethane	ND	10	856 J	5000	ND	5000	ND	2500
cis/trans-1,2-Dichloroethene	ND	10	ND	5000	ND	5000	ND	2500
Chloroform	ND	10	ND	5000	ND	5000	ND	2500
1,2-Dichloroethane	ND	10	ND	5000	ND	5000	ND	2500
2-Butanone	ND	10	ND	5000	ND	5000	ND	2500
1,1,1,-Trichloroethane	ND	10	ND	5000	ND	5000	ND	2500
Carbon tetrachloride	ND	10	ND	5000	ND	5000	ND	2500
Vinyl Acetate	ND	10	ND	5000	ND	5000	ND	2500
Bromodichloromethane	ND	10	ND	5000	ND	5000	ND	2500
1,2-Dichloropropane	ND	10	ND	5000	ND	5000	ND	2500
Trichloroethene	ND	10	ND	5000	ND	5000	ND	2500
Benzene	ND	10	ND	5000	ND	5000	ND	2500
cis-1,3-Dichloropropene	ND	10	960 J	5000	ND	5000	ND	2500
Dibromochloromethane	ND	10	ND	5000	ND	5000	ND	2500
trans-1,3-Dichloropropene	ND	10	ND	5000	ND	5000	ND	2500
1,1,2-Trichloroethane	ND	10	ND	5000	ND	5000	ND	2500
Bromoform	ND	10	ND	5000	ND	5000	ND	2500
4-Methyl-2-Pentanone	3.1 J	10	ND	5000	ND	5000	ND	2500
2-Hexanone	ND	10	ND	5000	ND	5000	ND	2500
1,1,2,2-Tetrachioroethane	ND	10	ND	5000	ND	5000	ND	2500
Tetrachioroethene	ND	10	ND	5000	ND	5000	ND	2500
Toluene	1.6	10	1400 JB	5000	730 JB	5000	550 JB	2500
Chlorobenzene	ND	10	ND	5000	ND	5000	ND	2500
Ethyl benzene	ND	10	1000 J	5000	ND	5000	ND	2500
Styrene	ND	10	6100	5000	130000	5000	40000	2500
m&p-Xylene	1.9	J 10	1800 JB	5000	ND	5000	520 JB	2500
o-Xylene	ND	10	790 J	5000	ND	5000	ND	2500
Methyl-tert-butyl Ether	ND	10	ND	5000	ND	5000	ND	2500
Tert – Butyl alcohol	ND	10	ND	5000	ND	5000	ND	2500

2 2 35.3

Table 1.9 Results of the BNA in Water Analysis
WA # 0-098 Pier Drum

	SAMPLE # : FILE : LOCATION* : COLLECTED : EXTRACTED : ANALYZED : INJECTED : MATRIX : DIL. FACT.: UNITS :	WBLK011895 ^PD010 01/18/95 01/30/95 13:36 WATER 1.0 ug/L		CD00636 ^PD011 MW-DUP 01/11/95 01/18/95 01/30/95 14:31 WATER 1.0 ug/L		CD00638 ^PD012 N-Pit 01/11/95 01/18/95 01/30/95 15:26 WATER 1.1 ug/L		CD00641 ^PD013 MW-N 01/11/95 01/18/95 01/30/95 16:21 WATER 1.0 ug/L		CD00635 ^PD014 MW-S 01/11/95 01/18/95 01/30/95 17:16 WATER 1.0 Ug/L	
COMPOUND		CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL
Phenol		ND	10	ND	10	ND	11	ND	11	ND	10 .
bis(-2-Chloroethyl)Ether	ND	10	ND	10	ND	11	ND	11	ND	10
2-Chlorophenol		ND	10	ND	10	ND	11	ND	11	ND	10
1,3-Dichlorobenzen	e	ND	10	ND	10	ND	11	ND	11	ND	10
1,4-Dichlorobenzen	e	ND	10	ND	10	ND	11	ND	11	ND	10
Benzyl alcohol		ND	10	ND	10	ND	11	ND	11	ND	10
1,2-Dichlorobenzen	e	ND	10	ND	10	ND	11	ND	11	ND	10
2-Methylphenol		ND	10	ND	10	ND	11	ND	11	ND	10
bis(2-Chloroisopro	pyl)ether	ND	10	ND	10	ND	11	ND	11 -	ND	10
4-Methylphenol	, ,	ND	10	ND	10	ND	11	ND	11	ND	10
N-Nitroso-Di-n-pro	pylamine	ND	10	ND	10	ND	11	ND	11	ND	10
Hexachloroethane		ND	10	ND	10	ND	11	ND	11	ND	10
Nitrobenzene		ND	10	ND	10	ND	11	ND	11	ND	10
Isophorone		ND	10	ND	10	ND	11	ND	11	ND	10
2-Nitrophenol		ND	10	ND	10	ND	11	ND	11	ND	10
2,4-Dimethylphenol		ND	10	ND	10	ND	11	ND	11	ND	10
bis(2-Chloroethoxy)methane	ND	10	ND	10	ND	11	ND	11	ND	10
2,4-Dichlorophenol	•	ND	10	ND	10	ND	11	ND	11	ND	10
1,2,4-Trichloroben	zene	ND	10	ND	10	ND	11	ND	11	ND	10
Naphthalene		ND	10	ND	10	ND	11	ND	11	ND	10
4-Chloroaniline		ND	10	ND	10	ND	11	ND	11	ND	10
Hexachlorobutadien	e	ND	10	ND	10	ND	11	ND	11	ND	10
4-Chloro-3-methylp		ND	10	ND	10	ND	11	ND	11	ND	10
2-Methylnaphthalen		ND	10	ND	10	ND	11	ND	11	ND	10
Hexachlorocyclopen		ND	10	ND	10	ND	11	ND	11	ND	10
2,4,6-Trichlorophe		ND	10	ND	10	ND	11	ND	11	ND	10
2,4,5-Trichlorophe		ND	50	ND	52	ND	54	ND	53	ND	52
2-Chloronaphthalen		ND	10	ND	10	ND	11	ND	11	ND	10
2-Nitroaniline		ND	50	ND	52	ND	54	ND	53	ND	52
Dimethylphthalate		ND	10	ND	10	ND	11	ND	11	ND	10
Acenaphthylene		ND	10	ND	10	ND	11	ND	11	ND'	10
3-Nitroaniline	_	ND	50	ND	52	ND	54	ND	53	ND	52
Acenaphthene		ND	10	ND	10	ND	11	ND	11	ND	10
2,4-Dinitrophenol		ND	50	ND	52	ND	54	ND	53	ND	52

^{*} denotes that, at the request of the task leader, the locations have been added

Table 1.9 (Cont) Results of the BNA in Water Analysis WA # 0-098 Pier Drum $\,$

	SAMPLE # : FILE : LOCATION* : COLLECTED : EXTRACTED : ANALYZED : INJECTED : MATRIX : DIL. FACT.: UNITS :	WBLK011895 ^PD010 01/18/95 01/30/95 13:36 WATER 1.0 ug/L		CD00636 ^PD011 MW-DUP 01/11/95 01/18/95 01/30/95 14:31 WATER 1.0 ug/L		CD00638 ^PD012 N-Pit 01/11/95 01/18/95 01/30/95 15:26 WATER 1.1 ug/L		CD00641 ^PD013 MM-N 01/11/95 01/18/95 01/30/95 16:21 WATER 1.0 Ug/L		CD00635	
COMPOUND		CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL
4-Nitrophenol		ND	50	ND	52	ND	54	ND	53	ND	52
Dibenzofuran		ND	10	ND	10	ND	11	ND	11	ND	10
2,6-Dinitrotoluene	!	ND	10	ND	10	ND	11	ND	11	ND	10
2,4-Dinitrotoluene	!	ND	10	ND	10	ND	11	ND	11	ND	10
Diethylphthalate		ND	10	ND	10	ND	11	ND	11	ND	10
4-Chlorophenyl-phe	nylether	ND	10	ND	10	ND	11	ND	11	ND	10
Fluorene		ND	10	ND	10	ND	11	ND	11	ND	10
4-Nitroaniline		ND	50	ND	52	ND	54	ND	53	ND	52
4,6-Dinitro-2-meth		ND	50	ND	52	ND	54	ND	53	ND	52
N-Nitrosodiphenyla		ND	10	ND	10	ND	11	ND	11	ND	10
4-Bromophenyl-phen	ylether	ND	10	ND	10	ND	11	ND	11	ND	10
Hexachlorobenzene		ND	10	ND	10	ND	11	ND	11	ND	10
Pentachlorophenol		ND	50	ND	52	ND	54	ND	53	ND	52
Phenanthrene		ND	10	ND	10	ND	11	ND	11	ND	10
Anthracene		ND	10	ND	10	ND	11	ND	11	ND	10
Carbazole		ND	10	ND	10	ND	11	MD	11	ND	10
Di-n-butylphthalat	e	2(J)	10	1(J,B)	10	38(B)	11	50(B)	11	2(J,B)	10
Fluoranthene		ND	10	ND	10	ND	11	ND	11	ND	10
Pyrene		ND	10	ND	10	ND	11	ND	11	ND	10
Butylbenzylphthala		ND	10	ND	10	ND	11	2(1)	11	ND	10
3,3'-Dichlorobenzi		ND	50	ND	52	ND	54	ND	53	ND	52
Benzo(a)anthracene		ND	10	ND	10	ND	11	ND	11	ND	10
Bis(2-Ethylhexyl)p	hthalate	1(J)	10	ND	10	2(J,B)	11	ND	1.1	3(J,B)	10
Chrysene		ND	10	ND	10	ND	11	ND	11	ND	10
Di-n-octylphthalat	e	ND	10	ND	10	ND	11	ND	11	ND	10
Benzo(b)fluoranthe		ND	10	ND	10	ND	11	ND	11	ND	10
Benzo(k)fluoranthe	ne	ND	10	ND	10	ND	11	ND	11	ND	10
Benzo(a)pyrene		ND	10	ND	10	ND	11	ND	11	ND	10
Indeno(1,2,3-cd)py		ND	10	ND	10	ND	11	ND	11	ND	10
Dibenzo(a,h)anthra		ND	10	ND	10	ND	11	ND	11	ND	10
Benzo(g,h,i)peryle	ne	ND	10	ND	10	ND	11	ND	11	ND	10

^{*} denotes that, at the request of the task leader, the locations have been added

Table 1.9 (Cont) Results of the BNA in Water Analysis WA # 0-098 Pier Drum

	SAMPLE # : FILE : LOCATION* : COLLECTED : EXTRACTED : ANALYZED : INJECTED : MATRIX : DIL. FACT.: UNITS :	West 01/1' 01/18 01/30 18:1' WATER 1.0	^PD015 West Pit (WP 01/11/95 01/18/95 01/30/95 18:11 WATER		CD00640 ^PD016 Pit) Field Blank 01/11/95 01/18/95 01/30/95 19:05 WATER 1.6 ug/L		534 19 1/95 8/95 0/95 6	CD00637 ^PD032 MW-E 01/11/95 01/18/95 02/01/95 11:37 WATER 1.1 ug/L	
COMPOUND		CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL
Phenol		ND	10	ND	16	ND	11	ND	11
bis(-2-Chloroethyl)	Ether	ND	10	ND	16	ND	11	ND	11
2-Chlorophenol		ND	10	ND	16	ND	11	ND	11
1,3-Dichlorobenzene		ND	10	ND	16	ND	11	ND	11
1,4-Dichlorobenzene		ND	10	ND	16	ND	11	ND	11
Benzyl alcohol		ND	10	ND	16	ND	11	ND	11
1,2-Dichlorobenzene		ND	10	ND	16	ND	11	ND	11
2-Methylphenol	15.45	ND	10	ND	16	ND	11	ND	11
bis(2-Chloroisoprop	yl)ether	ND	10	ND	16	ND	11	ND	11
4-Methylphenol		ND	10	ND	16	ND	11	ND	11
N-Nitroso-Di-n-prop	ylamine	ND	10	ND	16	ND	11	ND	11
Hexachloroethane		ND	10	ND	16	ND	11	ND	11
Nitrobenzene		ND	10	ND	16	ND	11	ND	11
Isophorone		ND	10	ND	16	ND	11	ND	11
2-Nitrophenol		ND	10	ND	16	ND	11	ND	11
2,4-Dimethylphenol		ND	10	ND	16	ND	11	ND	11
bis(2-Chloroethoxy)	methane	ND	10	ND	16	ND	11	ND	11
2,4-Dichlorophenol		ND	10	ND	16	ND	11	ND	11
1,2,4-Trichlorobenzo	ene	ND	10	ND	16	ND	11	ND	11
Naphthalene		ND	10 10	ND	16 16	ND	11 11	ND	11 11
4-Chloroaniline		ND ND	10	ND ND	16	ND ND	11	ND ND	11
Hexachlorobutadiene	anal	ND	10	ND	16	ND	11	ND ND	11
4-Chloro-3-methylph	enot	ND	10	ND	16	ND	11	ND ND	11
2-Methylnaphthalene	ndiana	ND	10	ND	16	ND	11	ND ND	11
Hexachlorocyclopenta		ND	10	ND ND	16	ND	11	ND ND	11
2,4,6-Trichlorophene 2,4,5-Trichlorophene		ND	52	ND	78	ND	57	ND ND	55
2-Chloronaphthalene	J.	ND	10	ND	16	ND	11	ND	11
2-Nitroaniline		ND	52	ND	78	ND ND	57	ND	55
Dimethylphthalate		ND ND	10	ND	16	ND	11	ND ND	33 11
Acenaphthylene		ND ND	10	ND	16	ND ND	11	ND ND	11
3-Nitroaniline		ND	52	ND ND	78	ND ND	57	ND	55
Acenaphthene		ND	10	ND	16	ND	11	ND	11
2,4-Dinitrophenol		ND	52	ND	78	ND	57	ND	55
e, a billier opileriot		,,,,,	~				٠.	NU	,,

 $^{^{\}star}$ denotes that, at the request of the task leader, the locations have been added

Table 1.9 (Cont) Results of the BNA in Water Analysis
WA # 0-098 Pier Drum

FILE LOCATION* COLLECTED EXTRACTED AMALYZED INJECTED MATRIX DIL. FACT.	: ^PD015		CD00634 ^PD019 MW-W 01/11/95 01/18/95 01/30/95 21:46 WATER 1.1 ug/L		CD00637 ^PD032 MW-E 01/11/95 01/18/95 02/01/95 11:37 WATER 1.1 ug/L			
COMPOUND	CONC.	MDL	CONC.	MDL	CONC.	MOL	CONC.	MDL
4-Nitrophenol	ND	52	ND	78	ND	57	ND	55
Dibenzofuran	ND	10	ND	16	ND	11	ND	11
2,6-Dinitrotoluene	ND	10	ND	16	ND	11	ND	11
2,4-Dinitrotoluene	ND	10	ND	16	ND	11	ND	11
Diethylphthalate	ND	10	ND	16	ND	11	ND	11
4-Chlorophenyl-phenylether	ND	10	ND	16	ND	11	ND	11
Fluorene	ND	10	ND	16	ND	11	ND	11
4-Nitroaniline	ND	52	ND	78	ND	57	ND	55
4,6-Dinitro-2-methylphenol	ND	52	ND	78	ND	57	ND	55
N-Nitrosodiphenylamine	ND	10	ND	16	ND	11	ND	11
4-Bromophenyl-phenylether	ND	10	ND	16	ND	11	ND	11
Hexach Lorobenzene	ND	10	ND	16	ND	11	ND	11
Pentachlorophenol	ND	52	ND	78	ND	57	ND	55
Phenanthrene	ND	10	ND	16	ND	11	ND	11
Anthracene	ND	10	ND	16	ND	11	ND	11
Carbazole	ND	10	ND	16	ND	11	ND	11
Di-n-butylphthalate	47(B)	10	4(J,B)	16	66(B)	11	1(J,B)	11
Fluoranthene	ND	10	ND	16	ND	11	ND	11
Pyrene	ND	10	ND	16	ND	11	ND	11
Butylbenzylphthalate	4(J)	10	ND	16	ND	11	ND	11
3,31-Dichlorobenzidine	ND	52	ND	78	ND	57	ND	55
Benzo(a)anthracene	ND	10	ND	16	ND	11	ND	11
Bis(2-Ethylhexyl)phthalate	ND	10	ND	16	1(J,B)	11	ND	11
Chrysene	ND	10	ND	16	ND	11	ND	11
Di-n-octylphthalate	ND	10	ND	16	ND	11	ND	11
Benzo(b)fluoranthene	ND	10	ND	16	ND	11	ND	11
Benzo(k)fluoranthene	ND	10	ND	16	ND	11	ND	11
Benzo(a)pyrene	ND	10	ND	16	ND	11	ND	11
Indeno(1,2,3-cd)pyrene	ND	10	ND	16	ND	11	ND	11
Dibenzo(a,h)anthracene	ND	10	ND	16	ND	11	ND	11
Benzo(g,h,i)perylene	ND	10	ND	16	ND	11	ND	11

^{*} denotes that, at the request of the task leader, the locations have been added

Table 1 / Ø	Perulis of Tental	ively Identi	fied Compounds	FOR BNA	in	Whiter
WA#	Results of Tentar	Dier	Drum			

Sample Number Lab File #

>PD 210

Unit Conversion Factor 100 L

CAS#	Compound Name	e ·	Q	RT	Conc.*
1.	Carboxylie Accid	aften 0.	-	31.02	11
2	Alkane	Culture		33.14	77
3.					
4.					
5.					
6.					
7.					
8.					
۹.					
10.					
11.		. ·			
12					<u> </u>
13.					
14.					
15.		·			
16.		·			<u> </u>
17.		·	<u> </u>		
18.					
19.					
20.		· · · · · · · · · · · · · · · · · · ·			
21.		···			
22		·	<u> </u>		
23.				1	

^{*} Estimated Concentration (Response Factor = 1.0)

8 3 mg

Table 1. 10 (cont.) Results of Tentatively Identified Compounds for BNA in Wafer WA# 0 - 098 Peer Drum

Sample Number Lab File # CD00616 >PD011

Unit Conversion Factor ers/L

CAS#	Compound Name	Q	RT	Conc.
1. 10544500	Sulfur, mel. St.		2783	15
2 -	Sulfur, mel. 58. Carboxylic Accid an Heno.	-	71.03	9
3.	J			
4.				
5.				
б.				
7.				
8.				
۷.				
10.				
11.				
12		ļ		<u> </u>
13.		<u> </u>		
14.			<u> </u>	
15.		ļ		
16.			ļ	
17.			ļ	
18.				
19.				
20.				
21.		<u> </u>		
22.				
23.				

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1. 16 (cont.) Results of Tentatively Identified Compounds for BNA in Water

WA# 0 - 098 Pier Drum

Sample Number Lab File #

CD 0 0638 >PD 0 12

Unit Conversion Factor 1.08 1.08

CAS#	Compound Name	Q	RT	Conc.
1. —	Unknown	-	6.70	8
2. –		_	15.95	10
3. —		-	16.76	5
4		-	78.85	5
5		_	31.03	12
6	Alkane CuH=n+2	-	35.0!	5
-	J		76.01	5
3.				
4.				
10.				
11.				
12.				
13.				
14.				
15.				
16.	,			
17.				·
18.				
19.				
20.				
21.				-
22.				
ಚ.				

^{*} Estimated Concentration (Response Factor = 1.0)

Table	1./0 (CORL)	Results of	Tentatively Iden	tified Com	pounds for	BNA	in	Water
	WA#	0.098	Pier	Drum				

Sample Number Lab File # CD00641 >PD013

Unit Conversion Factor 98/L 1.65

CAS#	Compound Name	Q	RT	Conc.*
1. 10544500	Sulfur. Sa	51	27.82	6
2 -		,	788c	
3. <u> </u>	Carbonglie Acid Cuffenos	_	31.05	
4	Carbonglie Acid Cuffenos Allame	_	34-11	5
5		-	35.03	5
6. —		_	36.03	5
7.	·			
8.				
Q _.				
10.				
t1.				
12.				
13.				
14.				
15.				
16.	•			
17.	•			
18.				
19.				
20.				·
21.				
22		1		<u> </u>
23.				

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1./0 (cont.) Results of Tentatively Identified Compounds for BNA in Wafer WA# 0-098 Pier Drum

Sample Number Lab File #

CD00615 >PD014

Unit Conversion Factor

CAS#	Compound Name	Q	RT	Conc.*
l. —	Chikneyn	-	6.79	(1
= 105HT200	Carboxy lie Acid Caffunda	80	27.81	8
3. —	Sulfur. 58 Carboxy lie Acid Cuffulle	-	31.02	9
4.				
5.				
6.				
7.	·			
8.				
υ.		<u> </u>		
10.				
11.				
12.		<u> </u>		
13.				
14.				
15.				
16.	,			
17.		<u> </u>		
18.				<u></u>
19.				
20.				
21.				
22.				
3 .				

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1. $\frac{j_0}{WA\# 0 - 09\%}$ Tentat	ively Identified Compounds	s for BNA	in	Water
WA# 0 - 098	Pier Drum			

Sample Number Lab File #

>PD 0/5

Unit Conversion Factor 1.64

CAS#	Compound Name	Q	RT	Conc.
1. —	Unknown	_	28.01	6
2	Carboxy lie Accid Cullinos	_	31.06	8
3. –	Carboxy lic Accid Cuffen 02 Phthalele	_	٦٧.٧٤	6
4.	A(kane	_	33.16	•
5. —		_	34.12	7
6		_	35.0L	10
7. —	V	-	36.06	7
8.				
Ų.				
10.				
11.				
12.				
13	·			
14.				
15.				
16.	,			
17.				
18.				
19.				
20.				
21.				
22.				
3.		1		

^{*} Estimated Concerts tion (Response Factor = 1.0)

Table 1	10,1000	Desults of	Tentatively l	identifier	1 Compounds	for BNA	in	Water_
1 abic		0.00	Di	Pr D	Ly m		•	
	~~~	<u>v. v. 771</u>						

Sample Number Lab File # CDOOGGO >PDOID

Unit Conversion Factor

		7		
CAS#	Compound Name	Q	RT	Conc.*
l. –	Carbonylie Acid		31.05	16
2				
3.		7		
4.				
5.				
6.				
7.				
8.	·			
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.	,			
17.				
18.				
19.				
20.				
21.				
<u> </u>				
ಚ.				

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1.70 (cont.) Results of Tentatively Identified Compounds for BNA in Water

WA# 0 - 098 Peer Drum

Sample Number Lab File # CD00634 >PD019

Unit Conversion Factor

CAS#	Compound Name	Q	RT	Conc.*
1	Unkneun	_	6.70	8
2 —			6.79	10
3. —		-	>1.28	7
4. 10544500	Sulfar St	18	27.87	18
5	Carboxylic Acid Cattente	-	28.86	7~
6.	cinknavn	-	28 26	7
7	Carboxylic Acid allands	<u> </u>	31.05	17
8. 🚤	A(kane aff m+2	_	32.18	7
9.		<u> -</u>	33.17	12
10. –		-	34.13	, '
11. —		<u> -</u>	35.06	100
12. –		<u> -</u>	36.07	17
13. —		-	37.23	11
14. —	V	_	38.57	7
15.				
16.	•			
17.	·			
18.				
19.				
20.				
21.				
77.				
23.		1		

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1. 10 (cont.) Results of Tentatively Identified Compounds for BNA in Water

WA# 0 - 098 Per Drum

Sample Number
Lab File #

>PD 032

Unit Conversion Factor

2.6.4	Compound Name	Q	RT	Conc.
CAS#	Benzose Acid Derivatives	+	13.65	9
l. –	Benzena ce fate	_	15.39	21
		ı	15-61	6
3. —	Chikuana	1	16.42	6
4	Ol Ala l'a anhadriale C8H4O3	95	16.66	7
5. 8 Jung	Phthalic anhydride C8H4O3	_	20.18	6
6. —	theknown	-	24.51	6
7	C OCurs S8	64	2791	21
x. (0244200	Swaar	1-	28.02	11
۹. ــ	Chenaun	1	28.22	8
10. —	Methyl Dricycloalkane Carboryli And C13 H20 O2	1		
11.		1_	28.3	7 12
12	Charles Acid after	+	31.0	
13. —	Carbonylic Acid altho	+-	35.4	
14. —	Alleme	+	133.0	1
15.		╁		-
16.	,	+		+-
17.		+		
18.		+		
19.				
20.		-		
21.		+		
22		+		
23.				

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1.11 Results of the BNA in Soil Analysis WA # 0-098 Pier Drum

		CONC.	MDL	CONC.	MOL	CONC.	MDL	CONC.	MD
UNITS	:	ug/K	(g	ug/l	(g	ug/l	(g	ug/K	(g
FINAL VOL	:	1		1		1		1	
AMT. USED	-	30		30		30		30	
% SOLID	:	100		81		83		82	
DIL. FACT.	:	1.0		1.0		1.0		1.0	
MATRIX	:	SOIL		SOIL		SOIL		SOIL	
INJECTED	:	12:14	•	13:0	3	14:04	•	15:00)
ANALYZED	:	01/31	/95	01/3	1/95	01/31	1/95	01/31	/95
EXTRACTED	:	01/13	1/95	01/13	3/95	01/13	3/95	01/13	5/9 5
COLLECTED	:	01/11	/95	01/1	1/95	01/1	1/95	01/11	/95
LOCATION*	:	~-		Sed :	S	Sed I	i		
FILE	:	^PD02	24	^PD02	25	^PD02	26	^PD02	27
SAMPLE #	:	SBLKC	11395	A0064	42	A0064	43	A0064	4

COMPOUND	CONC.	MDL	CONC.	MOL	CONC.	MDL	CONC.	MDL
Phenol	ND	330	ND	407	ND	398	ND	402
bis(-2-Chloroethyl)Ether	ND	330	ND	407	ND	398	ND	402
2-Chlorophenol	ND	330	ND	407	ND	398	ND	402
1,3-Dichlorobenzene	ND	330	ND	407	ND	398	ND	402
1.4-Dichlorobenzene	ND	330	ND	407	ND	398	ND	402
Benzyl alcohol	ND	330	ND	407	ND	398	ND	402
1,2-Dichlorobenzene	ND	330	ND	407	ND	398	ND	402
2-Methylphenol	ND	330	ND	407	ND	398	ND	402
bis(2-Chloroisopropyl)ether	ND	330	ND	407	ND	398	ND	402
4-Methylphenol	ND	330	ND	407	ND	398	ND	402
N-Nitroso-Di-n-propylamine	ND	330	ND	407	ND	398	ND	402
Hexachloroethane	ND	330	ND	407	ND	398	ND	402
Nitrobenzene	ND	330	ND	407	ND	398	ND	402
Isophorone	ND	330	ND	407	ND	398	ND	402
2-Nitrophenol	ND	330	ND	407	ND	398	ND	402
2.4-Dimethylphenol	ND	330	ND	407	ND	398	ND	402
bis(2-Chloroethoxy)methane	ND	330	ND	407	ND	398	ND	402
2.4-Dichlorophenol	ND	330	ND	407	ND	398	ND	402
1,2,4-Trichlorobenzene	ND	330	ND	407	ND	398	ND	402
Naphthalene	ND	330	ND	407	ND	398	ND	402
4-Chloroaniline	ND	330	ND	407	ND	398	ND	402
Hexach Lorobutadiene	ND	330	ND	407	ND	398	ND	402
4-Chloro-3-methylphenol	ND	330	ND	407	ND	398	ND	402
2-Methylnaphthalene	ND	330	ND	407	ND	398	ND	402
Hexachlorocyclopentadiene	ND	330	ND	407	ND	398	ND	402
2.4.6-Trichlorophenol	ND	330	ND	407	ND	398	ND	402
2,4,5-Trichlorophenol	ND	1650	ND	2037	ND	1988	ND	2012
2-Chloronaphthalene	ND	330	ND	407	ND	398	ND	402
2-Nitroaniline	ND	1650	ND	2037	ND	1988	ND	2012
Dimethylphthalate	ND	330	ND	407	ND	398	ND	402
Acenaphthylene	ND	330	ND	407	ND	398	ND	402
3-Nitroaniline	ND	1650	ND	2037	ND	1988	ND	2012
Acenaphthene	ND	330	ND	407	ND	398	ND	402
2.4-Dinitrophenol	ND	1650	ND	2037	ND	1988	ND	2012

f * denotes that, at the request of the task leader, the locations have been added

2 1 - 0332

Table 1.11 (Cont) Results of the BNA in Soil Analysis WA # 0-098 Pier Drum

SAMPLE #	:	SBLK011395	A00642	A00643	A00644
FILE	:	^PD024	^PD025	^PD026	^PD027
LOCATION*	:		Sed S	Sed N	
COLLECTED	:	01/11/95	01/11/95	01/11/95	01/11/95
EXTRACTED	:	01/13/95	01/13/95	01/13/95	01/13/95
ANALYZED	:	01/31/95	01/31/95	01/31/95	01/31/95
INJECTED	:	12:14	13:08	14:04	15:00
MATRIX	:	SOIL	SOIL	SOIL	SOIL
DIL. FACT	. :	1.0	1.0	1.0	1.0
% SOLID	:	100	81	83	82
AMT. USED	:	30	30	30	30
FINAL VOL	:	1	1	1	1
UNITS	:	ug/Kg	ug/Kg	ug/Kg	ug/Kg

COMPOUND	CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL
4-Nitrophenol	ND	1650	ND	2037	ND	1988	ND	2012
Dibenzofuran	ND	330	ND	407	ND	398	ND	402
2,6-Dinitrotoluene	ND	330	ND	407	ND	398	ND	402
2,4-Dinitrotoluene	ND	330	ND	407	ND	398	ND	402
Diethylphthalate	ND	330	87(J)	407	ND	398	ND	402
4-Chlorophenyl-phenylether	ND	330	ND	407	ND	398	ND	402
Fluorene	ND	330	ND	407	ND	398	ND	402
4-Nitroaniline	ND	1650	ND	2037	ND	1988	ND	2012
4,6-Dinitro-2-methylphenol	ND	1650	ND	2037	ND	1988	ND	2012
N-Nitrosodiphenylamine	ND	330	ND	407	ND	398	ND	402
4-Bromophenyl-phenylether	ND	330	ND	407	ND	398	ND	402
Hexach Lorobenzene	ND	330	ND	407	ND	398	ND	402
Pentachlorophenol	ND	1650	ND	2037	ND	1988	ND	2012
Phenanthrene	ND	330	ND	407	ND	398	ND	402
Anthracene	ND	330	ND	407	ND	398	ND	402
Carbazole	ND	330	ND	407	ND	398	ND	402
Di-n-butylphthalate	2400	330	2765(B)	407	3029(B)	398	1581(B)	402
Fluoranthene	ND	330	ND	407	ND	398	ND	402
Pyrene	ND	330	ND	407	ND	398	ND	402
Butylbenzylphthalate	150(J)	330	179(J,B) 407	156(J,8	398	ND	402
3,3'-Dichlorobenzidine	ND	1650	ND	2037	ND	1988	ND	2012
Benzo(a)anthracene	ND	330	ND	407	ND	398	ND	402
Bis(2-Ethylhexyl)phthalate	620	330	338(J,B) 407	ND	398	39(J,B	
Chrysene	ND	330	ND	407	ND	398	ND	402
Di-n-octylphthalate	ND	330	ND	407	ND	398	ND	402
Benzo(b)fluoranthene	ND	330	ND	407	ND	398	ND	402
Benzo(k)fluoranthene	ND	330	ND	407	ND	398	ND	402
Benzo(a)pyrene	ND	330	ND	407	ND	398	ND	402
Indeno(1,2,3-cd)pyrene	ND	330	ND	407	ND	398	ND	402
Dibenzo(a,h)anthracene	ND	330	ND	407	ND	398	ND	402
Benzo(g,h,i)perylene	ND	330	ND	407	ND	398	ND	402

^{*} denotes that, at the request of the task leader, the locations have been added

Table 1./2				RAIA :-	Soil
Table 1./	Results of Te	intatively identi	ned Compoun	ge tot Mall in	
WA#	0.048	Dier	Drum		

Sample Number Lab File # SPLKO1895

Unit Conversion Factor 98/kg

CAS#	Compound Name	·	Q	RT	Conc.
1. 119619	Benzophenone	CILHOD	89	1.93ء	1000
2 _	Chikuawa		1	22.84	410
3	Phthalate		-	25.10	23c
4. —	Alcohle	Cuffen+20	-	₹0.0€	170
5	Alkene/Cycloalkane	anthen	1	32.16	3306
6	Alkene/Cycloalkane Phosphonic Acid derivative		1	34.19	270
7.	Alcohle	Cu Hante O	_ /	18.81	30-
8. —	chikuova		1	<i>→8.48</i>	170
۷. —		·	1	40.37	170
10	l V		•	42.69	,
11.					
12.					
13.	•				
14.					
15.					
16.				1	
17.					
18.		برائند المال الفوالدوا فيوسي ميونون			,
19.					
20.					
21.		 			
22				1	
<u>-3</u> .					

^{*} Estimated Concentration (Pesnorise Postor = 1.0)

Table 1/2(cont.) Results of Tentatively Identified Compounds for BNA in Soil

WA# 0 - 098 Pier Drum

Sample Number Lab File #

A00642 >PD 025

Unit Conversion Factor 48/kg

CAS	S#	Compound Name			RT	Conc.*
l.	100414	Ethyl Benzone	Cafilo	97	6.52	580
2	85449	Phthalic anhydride	C2H403	99	16.65	780
3,	<u></u>	Cycloalkans	Cuffen	-	19.22	700
4.	119619	Benzophenone	CBH100	84	21.94	78e
5.		Cyclealkane/	Cuffen	_	22.37	250
6.		Phthalate		-	25.11	250
7.	-	Alkane	Cutlen +2	1	30.06	330
8.				1	31.13	210
4,		chehown		+	ع.رد	2700
10.	. ~	Ackane	Cublente	-	33.15	450
i 1.	_			Į	34.11	aro
12.	~			1	75.05	660
13.				-	36.03	580
14.	_		·	-	37.19	290
15.	-	J		_	38.5C	310
16.	_	Unterown.		-	39.10	660
i7.	-		·	-	40.1¢	290
18.	-			-	40.37	490
19.	-			-	41.69	540
20.						
21.						
22.						
23.						

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1.	12 (cont.)	Results of	Tentatively 1	identified	Compounds	for BNA	in .	<u> </u>
	WA#	U - U47		CF PI	ч			

Sample Number Lab File #

200643 20026 Unit Conversion Factor 40,10

CAS#	Compound Name	Compound Name		RT	Conc.*
1. 119619	Benzophenone	CHHIOD	64	21.93	840
-	Phthalate		_	25.10	320
3. 10544500	Sulfur	58	88	27.84	320
4	Carbonylie Acid		_	30.07	240
5. –	Chiknown		_	32.18	4400
6. —	j			32.80	240
7	Alcohle	Gultzute O	1	34.19	240
×. —	Chifmanin		_	35.05	360
۹. ــ	Alkane	allings	+	36.04	240
10	J		-	38.53	2
11	Alcohle	a Hante O	_	38.8€	362
12 —	chepranen			₹.99	284
13. —			_	39.11	560
14. —			_	40.37	40
15. —	\		_	42.69	640
16.		•			
17.		·			·
18.					
19.					
20.					
21.					
72					
23.					

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1/2 (cont.) Results of Tentatively Identified Compounds for BNA in Social WA# 0 - 098 Poer Drum

Sample Number Lab File # A00644 >PD 027

Unit Conversion Factor 40.6

CAS#	Compound Name	Q	RT	Conc.
1. 119619	Benjorhenene C13H100	75	21.95	2000
2 -	Chenoun	-	لالا.بد	200
3. –	unperoun Phthalate	-	25.10	280
4	Unknown	_	30.06	200
5		_	72.16	3800
6. –		<u> -</u>	34.18	?ec
7. –		<u> -</u>	35.ed	205
8. —		<u> </u>	38.77	40
4		-	38.97	240
10. —		-	39.10	490
11		<u> -</u>	4037	240
12 _		<u> </u> -	42.67	570
13. —	l v	_	43.11	530
14.				
15.				
16.	•			
17.				·
18.				
19.				
20.				
21.				
22				
23.				

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1.13 Results of the TCLP BNA Analysis
WA # 0-098 Pier Drum

SAMPLE # : WBLK011995 B00525 B00526 BE00527 ^PD003 A00643 ^PD004 A00644 FILE ^PD001 ^PD002 A00642 LOCATION : 01/18/95 01/19/95 COLLECTED : 01/18/95 01/18/95 EXTRACTED : 01/19/95 01/19/95 01/19/95 ANALYZED : INJECTED : 01/27/95 01/27/95 01/27/95 01/27/95 18:59 19:56 20:54 21:52

COMPOUND CONC. MDL CON		MATRIX : DIL. FACT.:		WATER WATER 1.0 1.3			WATEI	Ř	WATER 1.6	
Phenol ND 10 ND 13 ND 11 ND 16	UNI	TS :	ug/	L	ug/l	L	ug/i		ug/l	-
Dis(-2-Chloroethyl)Ether	COMPOUND		CONC.	MDL	CONC.	MDL	CONC.	MDL	CONC.	MDL
2-Chlorophenol	Phenol		ND	10	ND	13	ND	11	ND	16
1,3-Dichlorobenzene	bis(-2-Chloroethyl)Eth	ier	ND	10	ND	13	ND	11	ND	16
1,4-Dichlorobenzene	2-Chlorophenol		ND	10	ND	13	ND	11	ND	16
Benzyl alcohol	1,3-Dichlorobenzene		ND	10	ND	13	ND	11	ND	16
1,2-Dichlorobenzene ND 10 ND 13 ND 11 ND 16 2-Methylphenol ND 10 ND 13 ND 11 ND 16 bis(2-Chloroisopropyl)ether ND 10 ND 13 ND 11 ND 16 4-Methylphenol ND 10 ND 13 ND 11 ND 16 N-Nitroso-Di-n-propylamine ND 10 ND 13 ND 11 ND 16 N-Nitroso-Di-n-propylamine ND 10 ND 13 ND 11 ND 16 N-Nitroso-Di-n-propylamine ND 10 ND 13 ND 11 ND 16 Nitrobenzene ND 10 ND 13 ND 11 ND 16 Nitrobenzene ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16	1,4-Dichlorobenzene		ND	10	ND	13	ND	11	ND	16
2-Methylphenol	Benzyl alcohol		ND	10	16	13	17	11	15(J)	16
bis(2-Chloroisopropyl)ether ND 10 ND 13 ND 11 ND 16 4-Methylphenol ND 10 ND 13 ND 11 ND 16 Nitroso-Di-n-propylamine ND 10 ND 13 ND 11 ND 16 Hexachloroethane ND 10 ND 13 ND 11 ND 16 Nitrobenzene ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 2-4-Dimethylphenol ND 10 ND 13 ND 11 ND 16 2-4-Dimethylphenol ND 10 ND 13 ND 11 ND 16 2-4-Direchlorophenol ND 10 ND 13<	1,2-Dichlorobenzene		ND	10	ND	13	ND	11	ND	16
4-Methylphenol ND 10 ND 13 ND 11 ND 16 N-Nitroso-Di-n-propylamine ND 10 ND 13 ND 11 ND 16 Nexachloroethane ND 10 ND 13 ND 11 ND 16 Nitrobenzene ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorone ND 10 ND 13 ND 11 ND 16 Isophorophorol ND 10 ND 13 ND 11 ND 16 Isophorophorol ND 10 ND 13 ND 11 ND 16 Isophorophorol ND 10 ND 13 ND 11 ND 16 Isophorophorol ND 10 ND 13 ND 11 ND 16 Isophorophorophorophorophorophorophoropho	2-Methylphenol		ND	10	ND	13	ND	11	ND	16
N-Mitroso-Di-n-propylamine ND 10 ND 13 ND 11 ND 16 Hexachloroethane ND 10 ND 13 ND 11 ND 16 Nitrobenzene ND 10 ND 13 ND 11 ND 16 Lsophorone ND 10 ND 13 ND 11 ND 16 Lsophorone ND 10 ND 13 ND 11 ND 16 2-Nitrophenol ND 10 ND 13 ND 11 ND 16 2-Alitrophenol ND 10 ND 13 ND 11 ND 16 2-4-Dimethylphenol ND 10 ND 13 ND 11 ND 16 2-4-Dichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Dichlorophenol ND 10 ND 13 ND 11 ND 16 1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 Naphthalene ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 Hexachlorobutadiene ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-4-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 Acenaphthylene ND 50 ND 63 ND 54 ND 82 Acenaphthylene ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16	bis(2-Chloroisopropyl)	ether	ND	10	ND	13	ND	11	ND	16
N-Nitroso-Di-n-propylamine	4-Methylphenol		ND	10	ND	13	ND	11	ND	16
Nexachloroethane		mine	ND	10	ND	13	ND	11	ND	16
Isophorone			ND	10	ND	13	ND	11	ND	16
2-Nitrophenol ND 10 ND 13 ND 11 ND 16 2,4-Dimethylphenol ND 10 ND 13 ND 11 ND 16 bis(2-Chloroethoxy)methane ND 10 ND 13 ND 11 ND 16 2,4-Dichlorophenol ND 10 ND 13 ND 11 ND 16 1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 Naphthalene ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 4-chtorocyclopentadiene ND 10 ND 13 ND 11 ND 16 2,4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16	Nitrobenzene		ND	10	ND	13	ND	11	ND	16
2-Nitrophenol ND 10 ND 13 ND 11 ND 16 2,4-Dimethylphenol ND 10 ND 13 ND 11 ND 16 bis(2-Chloroethoxy)methane ND 10 ND 13 ND 11 ND 16 2,4-Dichlorophenol ND 10 ND 13 ND 11 ND 16 1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 Naphthalene ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 Hexachlorobutadiene ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2,4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16	Isophorone		ND	10	ND	13	ND	11	ND	16
2,4-Dimethylphenol ND 10 ND 13 ND 11 ND 16 bis(2-Chloroethoxy)methane ND 10 ND 13 ND 11 ND 16 2,4-Dichlorophenol ND 10 ND 13 ND 11 ND 16 1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 Naphthalene ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-4,6-Trichlorophenol			ND	10	ND	13	ND	11	ND	
bis(2-Chloroethoxy)methane NO 10 ND 13 ND 11 ND 16 2,4-Dichlorophenol ND 10 ND 13 ND 11 ND 16 1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 Naphthalene ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 Reachlorocyclopentadiene ND 10 ND			ND	. 10	ND	13	ND	11	ND	16
2,4-Dichlorophenol ND 10 ND 13 ND 11 ND 16 1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 Naphthalene ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 4-Chloroanethylphenol ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 4-A,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND		hane	ND		ND	13	ND			
1,2,4-Trichlorobenzene ND 10 ND 13 ND 11 ND 16 Naphthalene ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 Hexachlorobutadiene ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 Hexachlorocyclopentadiene ND 10 ND 13 ND 11 ND 16 2-4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND			ND	10	ND	13	ND	11	ND	16
Naphthalene ND 10 ND 13 ND 11 ND 16 4-Chloroaniline ND 10 ND 13 ND 11 ND 16 Hexachlorobutadiene ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2-Chloronaphthalene ND 10 ND	- •	•	ND	10	ND	13	ND	11	ND	
4-Chloroaniline ND 10 ND 13 ND 11 ND 16 Hexachlorobutadiene ND 10 ND 13 ND 11 ND 16 4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 Hexachlorocyclopentadiene ND 10 ND 13 ND 11 ND 16 2-4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate <	• •		ND	10	ND	13	ND	11	ND	16
4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 Hexachlorocyclopentadiene ND 10 ND 13 ND 11 ND 16 2,4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate ND 10 ND 13 ND 11 ND 16 Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND	•		ND	10	ND	13	ND	11	ND	16
4-Chloro-3-methylphenol ND 10 ND 13 ND 11 ND 16 2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 Hexachlorocyclopentadiene ND 10 ND 13 ND 11 ND 16 2,4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate ND 10 ND 13 ND 11 ND 16 Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND	Hexach Lorobutadiene		ND	10	ND	13	ND	11	ND	16
2-Methylnaphthalene ND 10 ND 13 ND 11 ND 16 Hexachlorocyclopentadiene ND 10 ND 13 ND 11 ND 16 2,4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16		ot	ND	10	ND	13	ND	11	ND	16
Hexachlorocyclopentadiene ND 10 ND 13 ND 11 ND 16 2,4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate ND 10 ND 13 ND 11 ND 16 Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16	2-Methylnaphthalene		ND	10	ND	13	ND	11	ND	16
2,4,6-Trichlorophenol ND 10 ND 13 ND 11 ND 16 2,4,5-Trichlorophenol ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate ND 10 ND 13 ND 11 ND 16 Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16		ene	ND	10	ND	13	ND	11	ND	16
2,4,5-Trichlorophenol ND 50 ND 63 ND 54 ND 82 2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate ND 10 ND 13 ND 11 ND 16 Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16			ND	10	ND	13	ND	11	ND	16
2-Chloronaphthalene ND 10 ND 13 ND 11 ND 16 2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate ND 10 ND 13 ND 11 ND 16 Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16			ND	50	ND	63	ND	54	ND	82
2-Nitroaniline ND 50 ND 63 ND 54 ND 82 Dimethylphthalate ND 10 ND 13 ND 11 ND 16 Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16			ND	10	ND	13	ND	11	ND	16
Dimethylphthalate ND 10 ND 13 ND 11 ND 16 Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16			ND	50	ND	63	ND	54	ND	82
Acenaphthylene ND 10 ND 13 ND 11 ND 16 3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16	Dimethylphthalate		ND	10	ND	13	ND	11	ND	
3-Nitroaniline ND 50 ND 63 ND 54 ND 82 Acenaphthene ND 10 ND 13 ND 11 ND 16	• •		ND	10	ND	13	ND	11		
Acenaphthene ND 10 ND 13 ND 11 ND 16			ND	50		63				
· · · · · · · · · · · · · · · · · · ·			ND	10	ND		ND	11		
	•		-							

Table 1.13 (Cont) Results of the TCLP BNA Analysis 2

	SAMPLE # FILE LOCATION COLLECTED EXTRACTED ANALYZED INJECTED MATRIX DIL. FACT	: : :	WBLKC ^PD00 01/19 01/27 18:59 WATER 1.0	7/95 7/95	B00529 ^PD000 A00647 01/18/ 01/27/ 19:56 WATER 1.3	95 195	B00526 ^PD003 A00643 01/18, 01/27, 20:54 WATER 1.1	5 195 195	BE005; ^PD004 A00644 01/19, 01/27, 21:52 WATER 1.6	795 795
	UNITS	:	ug/L		ug/L		ug/L		ug/L	
COMPOUND			CONC.	MOL	CONC.	MOL	CONC.	HDL	CONC.	MDL
4-Nitrophenol			ND	50	ND	63	ND	54	ND	82
Dibenzofuran			ND	10	ND	13	ND	11	ND	16
2,6-Dinitrotoluene	•		ND	10	ND	13	ND	11	ND	16
2,4-Dinitrotoluene			ND	10	ND	13	ND	11	ND	16
Diethylphthalate			ND	10	5(J)	13	2(J)	11	ND	16
4-Chiorophenyl-phe	nylether		ND	10	ND	13	ND	11	ND	16
Fluorene			ND	10	ND	13	ND	11	ND	16
4-Nitroaniline			ND	50	ND	63	ND	54	ND	82 .
4,6-Dinitro-2-meth	ylphenol		ND	50	ND	63	ND	54	ND	82 4
N-Nitrosodiphenyla	mine		ND	10	ND	13	ND	11	ND	16
4-Bromophenyl-pher	ylether		ND	10	ND	13	ND	11	ND	16
Hexachlorobenzene			ND	10	ND	13	ND	11	ND	16
Pentachlorophenol			ND	50	ND	63	ND	54	ND	82
Phenanthrene			ND	10	ND	13	ND	11	ND	16
Anthracene			ND	10	ND	13	ND	11	ND	16
Carbazole			ND	10	ND	13	ND	11	ND	16
Di-n-butylphthalat	e		46	10	56(B)	13	45(B)	11	91(B)	16
Fluoranthene			ND	10	ND	13	ND	11	ND	16
Pyrene			ND	10	ND	13	ND	11	ND	16
Butylbenzylphthala	te		5(J)	10	ND	13	4(J,8)	11	2(J,B)	16
3,31-Dichlorobenzi	dine		ND	50	ND	63	ND	54	ND	82
Benzo(a)anthracene			ND	10	ND	13	ND	11	ND	16
Bis(2-Ethylhexyl)p	hthalate		11	10	12(J,B)	13	11(B)	11	15(J,B)	16
Chrysene			ND	10	ND	13	ND	11	ND	16
Di-n-octylphthalat	e		ND	10	ND	13	ND	11	ND	16
Benzo(b)fluoranthe	ne		ND	10	ND	13	ND	11	ND	16
Benzo(k)fluoranthe	ne		ND	10	ND	13	ND	11	ND	16
Benzo(a)pyrene			ND	10	ND	13	ND .	11	ND	16
Indeno(1,2,3-cd)py	rene		ND	10	ND	13	ND	11	ND	16
Dibenzo(a,h)anthra	cene		ND	10	ND	13	ND	11	ND	16
<pre>Benzo(g,h,i)peryle</pre>			ND	10	ND	13	ND	11	ND	16

Table 1. <u>//</u> WA#	Results of Te	entatively ident	ified Compou	nds for	BNA	in TCLY	
WA#	<u>v. 047</u>	PCPF	Drum				

Sample Number Lab File # >PD 801

Unit Conversion Factor 18/L

CAS#	Compound Name	Q	RT	Conc.
1. —	Ethylhexyl diphenyl Phosphoric Acid C 20 Hz704	_	31.04	5
2 _	Ethylheryl diphenyl Phosphoric Acid	-	31.73	8
3.	C 20 H 2704	0_		
4.				
3. ·				
6.				
7.				
8.				
۷.				
10.				
11.				
12.				
13.				
14.				
15.				
16.	·			
17.	·			
18.				
19.				
20.				
21.				
22				
23.				

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1.14 (cont.) Results of Tentatively Identified Compounds for BNA in T(LP) WA# 0-098 Peer Drum

Sample Number
Lab File #

B00525 >PD002

Unit Conversion Factor 135

CAS#	Compound Name	Q	RT	Conc.
l. —	llyknown	_	670	B
2 _			15.21	16
3.	Phonyl Butanone Ciotal	-	15.21	11
4. —	Carlos Acid	ł	17.27	(0
5. —	Alkene / Cycloalkane CuHzn Phenel, 4-1-methyl-1-phonyletryl) Cystuso	_	19.14	16
6. 599644	Phenal 4-11-methyl-1-phonyletryl) Gotto0	99	25.13	6
7. –	Carbonylic Acid	-	31.05	8
з.	J			
۹.				
10.				
11.				
12.				
13.				
14.				
l 5 .	t of the state of			
16.	,			
17.				
18.				
19.				
20.				
21.				
22				
3.				

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1.14 (cont.) Results of Tentatively Identified Compounds for BNA in TCLP. WA# 0 - 098 Per Drum

2

Sample Number

Lab File #

Lab

CAS#	Compound Name	Q	RT	Conc
1. —	Chikuma	-	668	٢
2 -		-	12.61	10
3			12.68	9
4. —		_	1ª.46	5
5. —		_	12.28	
6		<u> -</u>	35.04	5
7.				
Х.				
4.				
10.				_
11.				
12.				
13.				
14.				
15.		<u> </u>		
16.	•			
17.	•			
18.				
19.				
20 .				
21.				
<u> </u>				
23.				

^{*} Estimated Concentration (Regionse Factor = 1.0)

Table 1.14 (cont.) Results of Tentatively Identified Compounds for BNA in TCLP WA# 0 - 098 Peer Drum

Sample Number
Lab File #

BE 00527 >PD004

Unit Conversion Factor 1.64

	·	_		<u></u>
CAS#	Compound Name	Q	RT	Conc
l. —	Unknown		71.03	8
2				
3.			-	
4.				
5.		<u> </u>		
6.	<i>!</i>			
- .	· · · · · · · · · · · · · · · · · · ·			
3.	·			
۷.				
10.				
11.				
12.				
13.				
14.				
l 5.	,			
16.	•			
17.	•			
1 8 .				
19.				
20.				
21.				
22				
3 .				

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1.15 Results of the BNA in Drum Liquid Analysis WA # 0-098 Pier Drum

	SAMPLE # : FILE : LOCATION : COLLECTED : EXTRACTED : ANALYZED : INJECTED : MATRIX : DIL. FACT.: UNITS :	WBLK07 -PD05* 02/14, 02/24, 13:18 WATER 1.0 ug/L	1 / 9 5	00912 ^PD054 REAC #9 01/10/95 02/14/95 02/24/95 17:25 DRUM LIQUID 595.0 ug/L		00913 ^PD055 REAC #2 01/10/95 02/14/95 02/24/95 18:38 DRUM LIQUID 435.0 ug/L		00914 ^PD056 REAC #4 01/10/95 02/14/95 02/24/95 19:55 DRUM LIQUID 4.0 ug/L CONC. MDL	
COMPOUND				CONC.	MDL	CONC.	MDL		
Phenol		ND	10	2186(J)	5950	4354	4350	430	40
bis(-2-Chloroethyl))Ether	ND	10	ND	5950	ND	4350	ND	40
2-Chlorophenol		ND	10	ND	5950	ND	4350	ND	40
1,3-Dichlorobenzen		ND	10	ND	5950	ND	4350	ND	40
1,4-Dichlorobenzene	e	ND	10	ND	5950	ND	4350	ND	40
Benzyl alcohol		ND	10	4712(J)	5950	1819(J)	4350	3300	40
1,2-Dichlorobenzen	e	ND	10	ND	5950	ND	4350	ND	40
2-Methylphenol		ND	10	ND	5950 5050	ND	4350	ND	40
bis(2-Chloroisopro	pyl)etner	ND	10 10	ND	5950 5950	ND ND	4350 4350	ND	40 40
4-Methylphenol		ND ND	10	ND ND	5950 5950	ND ND	4350 4350	24(J) ND	40
N-Nitroso-Di-n-prop Hexachloroethane	pytamine	ND ND	10	ND	5950	ND	4350	ND ND	40
Nitrobenzene		ND	10	ND	5950	ND	4350	ND	40
Isophorone		ND	10	ND	5950	ND	4350	ND	40
2-Nitrophenol		ND	10	ND	5950	ND	4350	ND	40
2,4-Dimethylphenol		ND	10	ND	5950	ND	4350	ND	40
bis(2-Chloroethoxy)	\methane	ND	10	ND	5950	ND	4350	ND	40
2,4-Dichlorophenol	, me charle	ND	10	ND	5950	ND	4350	ND	40
1.2.4-Trichlorobenz	zene	ND	10	ND	5950	ND	4350	ND	40
Naphthalene		ND	10	ND	5950	ND	4350	ND	40
4-Chloroaniline		ND	10	ND	5950	ND	4350	ND	40
Hexachlorobutadiene	2	ND	10	ND	5950	ND	4350	ND	40
4-Chloro-3-methylph	nenol	ND	10	ND	5950	ND	4350	ND	40
2-Methylnaphthalene	e	ND	10	ND	5950	ND	4350	ND	40
Hexachlorocyclopent		ND	10	ND	5950	ND	4350	ND	40
2,4,6-Trichloropher		ND	10	ND	5950	ND	4350	ND	40
2,4,5-Trichloropher		ND	50	ND	29750	ND	21750	ND	200
2-Chloronaphthalene	2	ND	10	ND	5950	ND	4350	ND	40
2-Nitroaniline		ND	50	ND	29750	ND	21750	ND	200
Dimethylphthalate		ND	10	638(J)	5950	1626(J)	4350	ND	40
Acenaphthylene		ND	10	ND	5950	ND	4350	ND	40
3-Nitroaniline		ND	50	ND	29750	ND	21750	ND	200
Acenaphthene		ND	10 50	ND	5950 29750	ND ND	4350 21750	ND ND	40 200
2,4-Dinitrophenol		ND	20	ND	29/30	NU	21/50	NU	200

2 1 9344

Table 1.15 (Cont) Results of the BNA in Drum Liquid Analysis WA # 0-098 Pier Drum

	FILE : ^PD051		01/ 02/ 02/ 17:	054 C #9 10/95 14/95 24/95 25 M LIQUID	01/ 02/ 02/ 18:	055 C #2 10/95 14/95 24/95 38 H LIQUID	4.0 ug/L		
COMPOUND		CONC.	MDL					CONC.	MDL
4-Nitrophenol		ND	50	ND	29750	ND	21750	ND	200
Dibenzofuran		ND	10	ND	5950	ND	4350	ND	40
2,6-Dinitrotoluene		ND	10	ND	5950	ND	4350	ND	40
2,4-Dinitrotoluene		ND	10	ND	5950	ND	4350	ND	40
Diethylphthalate		ND	10	ND	5950	ND	4350	ND	40
4-Chlorophenyl-phen	ylether	ND	10	ND	5950	ND	4350	ND	40
Fluorene		ND	10	ND	5950	ND	4350	ND	40
4-Nitroaniline		ND	50	ND	29750	ND	21750	ND	200
4,6-Dinitro-2-methy		ND	50	ND	29750	ND	21750	ND	200
N-Nitrosodiphenylam		ND	10	ND	5950	ND	4350	ND	40
4-Bromophenyl-pheny	lether	ND	10	ND	5950	ND	4350	ND	40
Hexachlorobenzene		ND	10	ND	5950	ND	4350	ND	40
Pentachlorophenol		ND	50	ND	29750	ND	21750	ND	200
Phenanthrene		ND	10	ND	5950	ND	4350	ND	40
Anthracene		ND	10	ND	5950	ND	4350	ND	40
Carbazole		ND	10	ND	5950	ND	4350	ND	40
Di-n-butylphthalate	!	ND	10	971(J)	5950	955(J)	4350	ND	40
Fluoranthene		ND	10	ND	5950	ND	4350	ND	40
Pyrene		ND	10	ND	5 9 50	ND	4350	ND	40
Butylbenzylphthalat		ND	10	1046(J)	5950	ND	4350	7(J)	40
3,3'-Dichlorobenzid	ine	ND	50	ND	29750	ND	21750	ND	200
Benzo(a)anthracene		ND	10	ND	5950	ND	4350	ND	40
Bis(2-Ethylhexyl)ph	thalate	ND	10	7123	5950	6423	4350	11(J)	40
Chrysene		ND	10	ND	5950	ND	4350	ND	40
Di-n-octylphthalate		ND	10	ND	5950	ND	4350	14(J)	40
Benzo(b)fluoranthen		ND	10	ND	5950	ND	4350	ND	40
Benzo(k)fluoranthen	e	ND	10	ND	5950	ND	4350	ND	40
Benzo(a)pyrene		ND	10	ND	5950	ND	4350	ND	40
Indeno(1,2,3-cd)pyr		ND	10	ND	5950	ND	4350	ND	40
Dibenzo(a,h)anthrac		ND	10	ND	5950	ND	4350	ND	40
Benzo(g,h,i)perylen	e _	ND	10	ND	5950	ND	4350	ND	40

Table 1.15 (Cont) Results of the BNA in Drum Liquid Analysis WA # 0-098 Pier Drum

	SAMPLE # : FILE : LOCATION : COLLECTED : EXTRACTED : ANALYZED : INJECTED : MATRIX : DIL. FACT.: UNITS :	^PD0 REAG 01/1 02/0 02/0 17:3	: #2 10/95 06/95 08/95 37 4 LIQUID
COMPOUND		CONC.	MDL
Phenol bis(-2-Chloroethyl 2-Chlorophenol 1,3-Dichlorobenzen 1,4-Dichlorobenzen Benzyl alcohol 1,2-Dichlorobenzen 2-Methylphenol bis(2-Chloroisopro 4-Methylphenol N-Nitroso-Di-n-pro Hexachloroethane Nitrobenzene Isophorone 2-Nitrophenol 2,4-Dimethylphenol bis(2-Chloroethoxy 2,4-Dichlorophenol 1,2,4-Trichloroben Naphthalene 4-Chloro-3-methylp 2-Methylnaphthalen Hexachlorocyclopen 2,4,6-Trichlorophe 2,4,5-Trichlorophe 2,4,5-Trichlorophe 2-Chloronaphthalen 2-Chloronaphthalen 2-Nitroaniline Dimethylphthalate Acenaphthylene 3-Nitroaniline	e e e pyl)ether pylamine)methane zene e henol e tadiene nol	ND ND ND ND ND ND ND ND ND ND ND ND ND N	500 500 500 500 500 500 500 500 500 500
Acenaphthene 2,4-Dinitrophenol		ND NO	500 2500

Table 1.15 (Cont) Results of the BNA in Drum Liquid Analysis WA # 0-098 Pier Drum

COMPOUND	SAMPLE # FILE LOCATION COLLECTED EXTRACTED ANALYZED INJECTED MATRIX DIL. FACT. UNITS	:	^PD04 REAC 01/10 02/00 02/00 17:37	#2 0/95 6/95 8/95 7 LIQUID
4-Nitrophenol			ND	2500
Dibenzofuran			ND	500
2,6-Dinitrotoluene			ND	500
2,4-Dinitrotoluene			ND	500
Diethylphthalate			ND	500
4-Chlorophenyl-pher	nylether		ND	500
Fluorene	•	1	ND	500
4-Nitroaniline		1	ND	2500
4,6-Dinitro-2-methy	/iphenoi	4	ND	2500
N-Nitrosodiphenylan	ni ne		ND	500
4-Bromophenyl-pheny	/lether	1	ND	500
Hexachiorobenzene		1	ND	500
Pentachlorophenol		1	ND	2500
Phenanthrene		1	ND	500
Anthracene		ı	ND	500
Carbazole		1	ND	500
Di-n-butylphthalate	•	1	ND	500
Fluoranthene			ND	500
Pyrene			ND	500
Butylbenzylphthalat			ND	500
3,3'-Dichlorobenzio	line		ND	2500
Benzo(a)anthracene		-	ND	500
Bis(2-Ethylhexyl)ph	ithalate		66(J)	500
Chrysene			ND	500
Di-n-octylphthalate			72(J)	500
Benzo(b)fluoranther			ND	500
Benzo(k)fluoranther	ne .		ND	500
Benzo(a)pyrene			ND	500
Indeno(1,2,3-cd)pyr			ND	500
Dibenzo(a,h)anthrac			ND	500
Benzo(g,h,i)peryler	ne e	,	ND	500

11:	.		g	CO RNA in	Orm Lizoud.
Table 1.10	Results of Ter	itativėjy identi	nea Compounas	tot Alait m	or the contract of the contrac
WA#	0.098	Pier	Drum		

Sample Number Lab File #

>PD051

Unit Conversion Factor ug/L

CAS#	Compound Name	Q	RT	Conc.*
1.	None famed			
2	i i			
3.				
4.				
5.		<u>.</u>	<u> </u>	<u> </u>
6.				
7.			ļ	
8.			<u> </u>	<u> </u>
9.			<u> </u>	<u> </u>
10.			ļ	
11.				<u> </u>
12.			<u> </u>	<u> </u>
13.			<u> </u>	
14.			<u> </u>	
15.			<u> </u>	
16.				
<u>17.</u>			 	
18.			ļ	<u> </u>
19.			 	
20.			 	
21.			 	
22.			-	
23.		ı	1	1

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1. 16(cont.) Results of Tentatively Identified Compounds for BNA in Drum Liquid WA# 0-098 Pier Drum

Sample Number Lab File #

>pdos8

Unit Conversion Factor ug/L

CAS#	Compound Name		Q	RT	Conc.
1. ~	Syrene or isomers	CoHo	-	7.30	710000
= 100527	Benzaldehyde	C>460	55	9.04	280000
3. —	Chikuswa		-	12.73	95000
4	Benzaic Acid or derivative			13.71	48000
5	Unknown	<i>:</i>		15.83	48000
6. —	J		_	15.94	89000
7. 8541119	Phthalic anhydride	C84403	99	16.70	190000
8	Buton- ane, - phonyl	CIOHIO	_	17.34	110000
9	Unknown		<u> </u>	17.95	31800C
10	Benzoic Aid derivatives		<u> -</u>	20.26	120000
11. —	Unknown			≥२.८५	30000
12	\		_	25.60	77000
13. —	Unknown Phthalete			10.85	65000
14				14,84	60000
15.	Kenknown			>9.40	95000
ا6. حـ	curknown Phthalate.		~	70.02	36000
17	<u> </u>		-	31.90	76000
18. —	Uhrknown		-	40.03	18000
19. —	J		-	40.12	24000
20	Unknown Phylicalete		_	43.71	30000
21.					
22.					
23.					

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1. 16 (cont.) Results of Tentatively Identified Compounds for BNA in Drum Liquid WA# 0. 098 Pier Drum

Sample Number Lab File #

00413 10X >Pd059

Unit Conversion Factor 4350

CAS#		Compound Name	Q	RT	Conc.
1	-	Styrene or isomers Ca. H8	_	7-37	5700000
2. /6	0527	Benzaldehijde C7460	67	9.07	\$60001
3.		Unknewn	_	12.77	(60000
4.		Carpoxylic Acid Capulos	_	13.04	44000
5.	_	Chiknews		14.29	14000
6.	_	d		14:43	17000
7. 8	25449	phthalic anhydride Coffees	96	16, 79	230000
3		Buten-one, phonyl Clotton	_	17.40	150001
9.		Chrown		17.95	21000
10.	_	Butanediane, methyl-phonyl- Citties		20.27	250
11.	_	unknown	_	23.29	61000
12.			_	23.49	120001
13.	-		-	2560	Boom
14	-			27.77	70006
15				29.46	52000
16	-	,	<u> -</u>	29.73	13 0000
17.			-	18.60	13000
18.	-		_	40.01	[(0000)
19.		V	-	40.12	10004
20.		theknown Phthalate	_	43.68	13000
21.					
<u> 22.</u>					
23.					

^{*} Estimated Concentration Response Factor = 1.0)

Table 1. 16(cont.) Results of Tentatively Identified Compounds for BNA in Drew Liquid
WA# 0 - 078 Pier Drum

Sample Number Lab File #

>Pd 060

Unit Conversion Factor 40.0

CAS	5#	Compound Name		Q	RT	Conc.
1.	_	Hyrene or isomers	C8H8	-	7.24	800
2	<i>;</i>	cinkuswa		_	7.47	640
3.	<u>.</u>	Carboxylic Acid	Culturos	1	17.5	23000
4.	_	Benzac Aad or derivation		_	1466	14000
5.	401812	Beuzene ace fate	GH8-01		15.64	1300
6.		cinference		_	15.97	3800
7.	<u> </u>	J J		_	16.11	280
ਤ.	501520	Benzene prepanoic Acid	C9 H 10 02	ex	17.06	440
9.		Dinettal Bensoic Acid issuers	COH, 002	_	17.26	240
10.		Chiknown			20.52	440
11.	-			-	20.70	280
12.				_	24.85	1300
13.					25.59	360
14.					28.0f	8800
15.		Carperylic Aud		_	28.48	6000
16.		unknown phthalate		_	28.46	640
17.		cheknown		-	28.88	600
18.				_	29.42	1600
19.	_			_	30.99	1200
20.		V		-	31.06	440
21.						
<u>:2.</u>						
ಚ.						

^{*} Estimated Concentration (Response Factor = 1.0)

Table 1. 16(cont.) Results of Tentatively Identified Compounds for BNA in Drew Ligard
WA# 0-098 Pier Drum

Sample Number Lab File #

00917 (Bottom) >Pd040

Unit Conversion Factor ug/6

CA	S#	Compound Name		Q	RT	Conc.*
1.	_	Signene or isomers (8			7.18	36000
2	100527	Benzaldehyde Cithe		58	9.05	1800
3.	_	Cheknenin		-	9.28	1790
4.		Ethyl Heranoic Acid Caffe	601	_	12.83	1600
5.	-	Unknown			14.30	500
6.		V	1.	_	14'47	1100
7.	85449	Phthalic anhydride co	4401	99	16.74	6500
ક.		Buten-One, phonyl C10 Unknown Benzene deprivationes	Hr. 0	-	17.43	300
۷.		Benzene de privatines			20.27	500
10.		chknown		_	≥3.24	}ι
11.					23.51	(60
12.				_	23.53	780
13.	<u></u>				25.63	480
14.					27-39	250
15.				_	29.53	250
16.	<u> </u>			_	29.72	350
17.				_	40.(2	501
18.	- -			_	40.21	550
19.	-			_	40.A	450
20.	<u> </u>	V		_	43.82	550
21.						
22.						~
23.						

^{*} Estimated Concentration (Response Factor = 1.

Table 1.17 Results of the BNA Analysis for the TCLP Extracts from the Drum Solids WA # 0 -098 Pier Drum

Cilent ID Location	Blank		00911 Composit 1,	8,9,11	00915 REAC #5		00916 Composit 3,7	7,10,13,14
	Conc	MDL	Conc	MDL	Conc	MDL	Conc	MDL
Compound	(ug/l)	(ug/l)	(ug/l)	(ug/l)	(ug/l)	(ug/l)	(ug/l)	(ug/l)
Pyridine	ND	10	ND	20	ND	20	ND	20
N-Nitrosodidimethylamine	ND	10	ND	20	ND	20	ND	20
Anlline	ND	10	ND	20	ND	20	ND	50
Bis(2-chloroethyl) ether 1,3-Dichlorobenzene	ND ND	10 10	ND ND	20 20	ND ND	20 20	ND ND	20 20
1.4 - Dichlorobenzene	ND	10	ND	20	ND	20	ND	20
1,2 - Dichlorobenzene	ND	10	ND	20	ND	20	ND	20
Benzvi alcohol	ND	10	1000	20	83	20	150	20
Bis(2-chlorolsopropyl) ether	ND	10	ND	20	ND	20	ND	20
-lexachioroethane	ND	10	ND	20	ND	20	ND	20
N-Nitrosodi-n-propylamine	ND	10	ND	20	ND	20	ND	20
Nitrobenzene	ND	10	ND	20	ND	20	ND	20
sophorone	ND	10	ND ND	20 20	ND ND	20	ND	20
Bis(2-chloroethoxy) methane 1,2,4-Trichlorobenzene	ND ND	10 10	ND	20	ND	20 20	ND ND	20 20
Naphthalene	ND	10	ND	20	ND	20	16 J	20
Benzoic Acid	ND	10	120	20	270	20	4800 E	20
- Chloroaniline	ND	10	ND	20	ND	20	ND	20
lexachlorobutadiene	ND	10	ND	20	ND	20	ND	20
? - Methylnaphthalene	ND	10	ND	20	ND	20	6 J	20
lexachlorocyclopentadiene	ND	10	ND	20	ND	20	ND	20
-Chloronaphthalene	ND	10	ND	20	ND	20	ND	20
-Nitroaniline	ND	10	ND	20	ND	20	ŅĐ	20
cenaphthylene	ND	10	ND	20	ND	20	ND	20
Dimethyl phthalate	ND	10	ND	20	ND	20	ND	20
2,6 - Dinitrotoluene	ND	10	ND ND	20 20	ND	20	ND	20
Acenaphthene I - Nitroaniline	ND ND	10 10	ND ND	20	ND ND	20 20	ND ND	20
Dibenzofuran	ND	10	ND	20	ND ND	20	ND	20 20
2.4 – Dinitrotoluene	ND	10	ND	20	ND	20	ND	20
luorene	ND	10	ND	20	ND	20	ND	20
Diethylphthalate	ND	10	ND	20	ND	20	ND	20
-Chlorophenyl phenyl ether	ND	10	ND	20	ND	20	ND	20
-Nitroaniline	ND	10	ND	20	ND	20	ND	20
N-Nitrosodiphenylamine	ND	10	ND	20	ND	20	ND	20
-Bromophenyl phenyl ether	ND	10	ND	20	ND	20	ND	20
lexachlorobenzene	ND	10	ND	20	ND	20	ND	20
henanthrene	ND	10 10	ND ND	20 20	ND ND	20	ND	20
Anthracene Di – n – butyl phthalate	ND ND	10	ND 40	20	ND 34	20 20	9 J ND	20 20
fluoranthene	ND	10	ND	20	ND	20	ND	20
Benzidine	ND	10	ND	20	ND	20	ND	20
Pyrene	ND	10	ND	20	ND	20	ND	20
Butyl benzyl phthalate	ND	10	ND	20	ND	20	ND	20
Benzo(a)anthracene	ND	10	ND	20	ND	20	ND	20
3' - Dichlorobenzidine	ND	10	ND	20	ND	20	ND	20
Chrysene	ND	10	ND	20	ND	20	ND	20
3is(2-ethylhexyl) phthalate	ND	10	5 J	20	ND	20	13 J	20
Di-n-octyl phthalate	ND	10	8 J	20	. 6 1		. B J	20
Benzo(b)fluoranthene	ND	10	ND	20	ND	20	ND	20
Benzo(k)fluoranthene	ND	10	ND ND	20 20	ND	20	ND	20
Benzo(a)pyrene ndeno(1,2,3 - c,d)pyrene	ND ND	10 10	ND	20	ND ND	20 20	ND ND	2(2)
Dibenzo(a,h)anthracene	ND	10	ND	20	ND	20	ND	2
Benzo(ghi)perylene	ND	10	ND	20	ND	20	ND	2
Phenol	ND	10	110	20	17 J		34	2
-Nitrophenol	ND	10	ND	20	ND	20	ND	2
2,4 - Dimethylphenol	ND	10	ND	20	ND	20	ND	2
- Chlorophenol	ND	10	ND	20	ND	20	ND	2
2.4 - Dichlorophenol	ND	10	ND	20	ND	20	ND	2
I – Chloro – 3 – methylphenol	ND	10	ND	20	ND	20	ND	2
2,4,6~Trichlorophenol	ND	10	ND	20	ND	20	ND	2
2.4 – Dinitrophenol	ND	10	ND	20	ND	20	ND	2
1,6 - Dinitro - 2 - methylphenol	ND	10	ND	20	ND	20	ND	2
-Nitrophenol	ND ND	10	ND ND	20	ND ND	20	ND	2
Pentachlorophenol	ND ND	10 10	ND ND	20 20	ND ND	20 20	ND ND	2
2,4,5-Trichlorophenol 2-Methylphenol	ND	10	860	20	ND ND	20	ND ND	2
				_0	,,,,	20		

Table 1.18 Results of the Pest/PCB Analysis for the Water Samples WA # 0-098 Pier Drum

Client ID Location*	WBLK111795		EF 00634 MW-W		EF 00635 MW-S		EF 00636 MW_DUP		EF 00637 MW-E		EF 00638 N Pit	
	CONC	MDL	CONC	MDL	CONC	MDL	CONC	MDL	CONC	MDL	CONC	MDL
Analyte	μg/L	μg/L	μg/L	µg/L	μ g/ L	μ g /L	μg/L	μg/L	μg/L	μg/L	μ g /L	μg /L
a-BHC	ND	0.02	ND	0.02	NĎ	0.02	ND	0.02	ND	0.02	ND	0.02
g~BHC	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	C.02	ND	0.02
ь-внс	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Heptachlor	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
d-BHC	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Aldrin	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Heptachlor Epoxide	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
g-Chlordane	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
a-Chlordane	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Endosulfan (l)	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
p,p'-D D E	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Dieldrin	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Endrin	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
p,p'-D D D	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Endosulfan (II)	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
p,p'-DDT	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Endrin Aldehyde	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Endosulfan Sulfate	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Methoxychlor	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Endrin Ketone	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02	ND	0.02
Toxaphene	ND	0.50	ND	0.54	ND	0.52	ND	0.52	ND	0.55	ND	0.54
Aroclor 1016	ND	0.25	ND	0.27	ND	0.26	ND	0.26	ND	0.27	ND	0.27
Aroclor 1221	ND	0.50	ND	0.54	ND	0.52	ND	0.52	ND	0.55	ND	0.54
Aroclor 1232	ND	0.25	ND	0.27	ND	0.26	ND	0.26	ND	0.27	ND	0.27
Aroclor 1242	ND	0.25	ND	0.27	ND	0.26	ND	0.26	ND	0.27	ND	0.27
Aroclor 1248	ND	0.25	ND	0.27	ND	0.26	ND	0.26	ND	0.27	ND	0.27
Aroclor 1254	ND	0.25	ND	0.27	ND	0.26	ND	0.26	ND	0.27	ND	0.27
Aroclor 1260	ND	0.25	ND	0.27	ND	0.26	ND	0.26	ND	0.27	ND	0.27

^{*} denotes that, at the request of the task leader, the locations have been added

Table 1.18 (Cont) Results of the Pest/PCB Analysis for the Water Samples WA # 0-098 Pier Drum

Client ID	EF 00		EF 00		EF OC	
Location*	West Pit		Field E		MW-	
	CONC	MDL	CONC	MDL	CONC	MDL
Analyte	μg/L	μg/L	μ g /L	µg/L	μg/L	μ g /L
a-BHC	ND	0.02	ND	0.03	ND	0.02
g~BHC	ND	0.02	ND	0.03	ND	0.02
b~BHC	ND	0.02	ND	0.03	ND	0.02
Heptachlor	ND	0.02	ND	0.03	ND	0.02
d-BHC	ND	0.02	ND	0.03	ND	0.02
Aldrin	ND	0.02	ND	0.03	ND	0.02
Heptachlor Epoxide	ND	0.02	ND	0.03	ND	0.02
g-Chlordane	ND	0.02	ND	0.03	ND	0.02
a-Chiordane	ND	0.02	ND	0.03	ND	0.02
Endosulfan (I)	ND	0.02	ND	0.03	ND	0.02
p,p'-DDE	ND	0.02	ND	0.03	ND	0.02
Dieldrin	ND	0.02	ND	0.03	ND	0.02
Endrin	ND	0.02	ND	0.03	ND	0.02
p,p'-D D D	ND	0.02	ND	0.03	ND	0.02
Endosultan (II)	ND	0.02	ND	0.03	ND	0.02
p,p'~D D T	ND	0.02	ND	0.03	ND	0.02
Endrin Aldehyde	ND	0.02	ND	0.03	ND	0.02
Endosulfan Sulfate	ND	0.02	ND	0.03	ND	0.02
Methoxychior	ND	0.02	ND	0.03	ND	0.02
Endrin Ketone	ND	0.02	ND	0.03	ND	0.02
Toxaphene	ND	0.52	ND	0.79	ND	0.54
Aroclor 1016	ND	0.26	ND	0.40	ND	0.27
Aroclor 1221	ND	0.52	ND	0.79	ND	0.54
Arocior 1232	ND	0.26	ND	0.40	ND	0.27
Aroclor 1242	ND	0.26	ND	0.40	ND	0.27
Aroclor 1248	ND	0.26	ND	0.40	ND	0.27
Aroclor 1254	ND	0.26	ND	0.40	ND	0.27
Aroclor 1260	ND	0.26	ND	0.40	ND	0.27

^{*} denotes that, at the request of the task leader, the locations have been added

Table 1.19 Results of the Pest/PCB Analysis for the Soil Samples WA # 0-098 Pier Drum
Based on Dry weight

Client ID	SBLK011395			A00643 A00644		SBLK	20795	A00642 Sed S		
Location *		. .		dN		d E		=_		
% Solid	11	ōo	8	13	ε	12	10	ŌO		31
		MDL		MDL		MDL		MDL		MDL
Analyte	μ g /Kg	μ g /Kg	μ g /Kg	μ g /Kg	μ g/ Kg	μg/Kg	μg/Kg	µg/Kg	μg/Kg	µg/Kg
a-BHC	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
g-BHC	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
b-BHC	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Heptachior	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
d-BHC	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Aldrin	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Heptachlor Epoxide	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
g-Chiordane	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
a-Chlordane	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Endosulfan (i)	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
p,p'-DDE	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	[~] 41
Dieldrin	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Endrin	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
p,p'-D D D	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Endosulfan (II)	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
p,p'-D D T	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Endrin Aldehyde	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Endosulfan Sulfate	ND	3.3	ND	4.0	ND	4.1	ND	.∕ 33	ND	41
Methoxychior	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Endrin Ketone	ND	3.3	ND	4.0	ND	4.1	ND	33	ND	41
Toxaphene	ND	83	ND	100	ND	100	ND	830	ND	1000
Aroclor 1016	ND	42	ND	50	ND	51	ND	420	ND	510
Aroclor 1221	ND	83	ND	100	ND	100	ND	830	ND	1000
Aroclor 1232	ND	42	ND	50	ND	51	ND	420	ND	510
Aroclor 1242	ND	42	ND	50	ND	51	ND	420	ND	510
Aroclor 1248	ND	42	ND	50	ND	51	ND	420	ND	510
Aroclor 1254	ND	42	ND	50	ND	51	ND	420	ND	510
Aroclor 1260	ND	42	ND	50	ND	51	ND	420	ND	510

^{*} at the request of the Task Leader, the actual locations have been added

Table 1.20 Results of the TCLP Pest/PCB Analysis WA # 0-098 Pier Drum

Client ID Location*	WBLK011995		-	C 00525 A 00642		0526 0643	CF 00527 A 00644		
	•	MDL		MDL		MDL		MDL	
Analyte	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L	μg/L	μ g /L	
a-BHC	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
g-BHC	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
b-BHC	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Heptachlor	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
d-BHC	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Aldrin	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Heptachlor Epoxide	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
g-Chiordane	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
a-Chlordane	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Endosulfan (l)	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
p,p'-DDE	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Dieldrin	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Endrin	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
p,p'-D D D	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Endosulfan (II)	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
p,p'-D D T	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Endrin Aldehyde	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Endosulfan Sulfate	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Methoxychlor	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Endrin Ketone	ND	0.02	ND	0.02	ND	0.02	ND	0.04	
Toxaphene	ND	0.50	ND	0.60	ND	0.55	ND	0.88	
Aroclor 1016	ND	0.25	ND	0.30	ND	0.27	ND	0.44	
Aroclor 1221	ND	0.50	ND	0.60	ND	0.55	ND	0.88	
Aroclor 1232	ND	0.25	ND	0.30	ND	0.27	ND	0.44	
Aroclor 1242	ND	0.25	ND	0.30	ND	0.27	ND	0.44	
Aroclor 1248	ND	0.25	ND	0.30	ND	0.27	ND	0.44	
Aroclor 1254	ND	0.25	ND	0.30	ND	0.27	ND	0.44	
Aroclor 1260	ND	0.25	ND	0.30	ND	0.27	ND	0.44	

^{*} denotes that, at the request of the task leader, the locations have been added

Table 1.21 Results of the Pest/PCB Analysis in Drum Liquid Samples WA # 0-098 Pier Drum

Client ID Location	WBLK0	20795	009 Rea	12 ic#9	00913 Top layer 00913 Bottom layer Reac#2 Reac#2			009 Res	14 ic#4	
An ab da	(1	MDL	=0	MDL		MDL	4.07	MDL	/1	MDL
Analyte	μg/L	μg/L	μg/L	μg/L	µg/L	µg/L	μg/L	μ g/L	μg/L	µg/L
a~BHC	ND	0.02	0.16	0.11	ND	0.11	ND	1000	ND	0.11
g~BHC	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
bBHC	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Heptachlor	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
d~BHC	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Aldrin	ND	0.02	ND	0.11	0.18	0.11	ND	1000	ND	0.11
Heptachlor Epoxide	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
g-Chlordane	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
a-Chiordane	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Endosulfan (I)	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
p,p'-D D E	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Dieldrin	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Endrin	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
p,p'-D D D	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Endosulfan (II)	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
p,p'-DDT	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Endrin Aldehyde	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Endosulfan Sulfate	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Methoxychlor	ND	0.02	ND	0.11	ND	0.11	ND	1000	ND	0.11
Endrin Ketone	ND	0.02	ND	0.11	ND	0.11	ND	⊭100 0	ND	0.11
Toxaphene	ND	0.50	ND	2.80	ND	2.80	ND	25000	ND	2.80
Aroclor 1016	ND	0.25	ND	1.40	ND	1.40	ND	13000	ND	1.40
Aroclor 1221	ND	0.50	ND	2.80	ND	2.80	ND	25000	ND	2.80
Aroclor 1232	ND	0.25	ND	1.40	ND	1.40	ND	13000	ND	1.40
Aroclor 1242	ND	0.25	ND	1.40	ND	1.40	ND	13000	ND	1.40
Aroclor 1248	ND	0.25	ND	1.40	ND	1.40	ND	13000	ND	1.40
Aroclor 1254	ND	0.25	ND	1.40	ND	1.40	ND	13000	ND	1.40
Aroclor 1260	ND	0.25	ND	1.40	ND	1.40	ND	13000	ND	1.40

Table 1.22 Results of the Pest/PCB Analysis in Drum Solid Samples WA # 0-098 Pier Drum

J 553

00911 00915 00916 Client ID REAC #5 Composit Composit Location* 1, 8, 9, 11 3, 7, 10, 13, 14 MDL Conc MDL MDL Conc MDL Conc Conc Analyte μg/L μg/L µg/L µg/L µg/L µq/L μg/L µa/L Aldrin ND 0.10 ND 0.18 ND 0.11 ND 0.25 a-BHC ND 0.10 ND 0.18 ND 0.11 ND 0.25 b-BHC ND 0.10 ND 0.18 ND 0.11 ND 0.25 g-BHC ND 0.10 ND 0.18 ND 0.11 ND 0.25 d-BHC ND ND ND 0.60 0.18 0.11 0.25 0.10 ND a-Chiordane ND 0.10 ND 0.18 0.11 ND 0.25 g-Chlordane ND 0.10 ND 0.18 ND 0.11 ND 0.25 ND ND 0.18 ND 0.11 ND p,p'-D D D 0.10 0.25 p,p'-D D E ND 0.10 ND 0.18 ND 0.11 ND 0.25 ND ND 0.11 p,p'-DDT ND 0.18 ND 0.25 0.10 Dieldrin ND 0.10 ND 0.18 ND 0.11 ND 0.25 ND 0.18 ND ND 0.11 Endosulfan (alpha) 0.10 1.0 0.25 Endosulfan (beta) ND 0.10 ND 0.18 ND 0.11 ND 0.25 Endosulfan Sulfate ND ND 0.10 0.18 ND 0.11 ND 0.25 ND ND 0.18 ND ND **Endrin** 0.10 0.11 0.25 Endrin Aldehyde ND 0.10 ND 0.18 ND 0.11 ND 0.25 ND ND 0.18 ND 0.11 ND **Endrin Ketone** 0.10 0.25 Heptachlor ND 0.10 ND 0.18 ND 0.11 ND 0.25 Heptachlor Epoxide ND 0.10 ND 0.18 ND 0.11 ND 0.25 ND ND Methoxychlor 0.10 0.18 ND 0.11 ND 0.25 ND Toxaphene ND 1.00 ND 1.8 1.1 ND 2.5 ND ND ND ND Arocior 1016 1.00 1.8 1.1 2.5 Aroclor 1221 ND ND ND ND 1.00 1.8 1.1 2.5 Aroclor 1232 ND 1.00 ND ND ND 1.8 1.1 2.5 Aroclor 1242 ND 1.00 ND 1.8 ND ND 2.5 1.1 ND ND 2.5 ND ND Aroclor 1248 1.00 1.8 1.1 Aroclor 1254 ND 1.00 ND 1.8 ND 1.1 ND 2.5 Aroclor 1260 ND 1.00 ND 1.8 ND ND 2.5 1.1

^{*} at the request of the Task Leader, the actual locations have been added

Table 1.23 Results of the Metals Analysis for the Water Samples WA # 0-098 Pier Drum

Client ID Location		Method Lai		A006 Unfilte		B006 Filter		A006 Unfilte		B006 Filter		A006 Unfilte	
Parameter	Analysis Method	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l
Aluminum	ICAP	ND	40	1200	40	1000	40	2300	40	870	40	16000	40
Antimony	AA-Fur	ND	5.6	ND	5.6	ND	5.6	ND	5.6	ND	5.6	ND	5.6
Arsenic	AA-Fur	ND	2.2	3.5	2.2	3.3	2.2	30	2.2	29	2.2	12	2.2
Barium	ICAP	ND	4.0	18	4.0	17	4.0	20	4.0	10	4.0	44	4.0
Beryllium	ICAP	ND	2.0	ND	2.0	ND	2.0	ND	2.0	ND	2.0	ND	2.0
Cadmium	AA-Fur	ND	0.22	ND	0.22	ND	0.22	ND	0.22	ND	0.22	ND	0.22
Calcium	ICAP	ND	100	4000	100	4000	100	17000	100	16000	100	2300	100
Chromium	AA-Fur	ND	2.8	ND	2.8	ND	2.8	8.6	2.8	4.5	2.8	16	2.8
Cobalt	ICAP	ND	6.0	ND	6.0	ND	6.0	ND	6.0	ND	6.0	ND	6.0
Copper	ICAP	ND	5.0	ND	5.0	ND	5.0	ND	5.0	ND	5.0	ND	5.0
Iron	ICAP	ND	10	9400	10	9100	10	950	10	560	10	4100	10
Lead	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2	3.0	2.2
Magnesium	ICAP	ND	500	4200	500	4100	500	6500	500	5800	500	1400	500
Manganese	ICAP	ND	2.0	4.0	2.0	3.0	2.0	18	2.0	13	2.0	5.0	2.0
Mercury	Cold Vapor	ND	0.20	ND	0.20	ND	0.20	ND	0.20	ND	0.20	ND	0.20
Nickel	ICAP	ND	10	ND	10	ND	10	ND	10	ND	10	ND	10
Potassium	ICAP	ND	2000	ND	2000	ND	2000	4900	2000	4400	2000	ND	2000
Selenium	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2	√ ND	2.2
Silver	AA -Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2
Sodium	ICAP	ND	500	29000	500	29000	500	25000	500	23000	500	3800	500
Thallium	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2
Vanadium	ICAP	ND	5.0	ND	5.0	ND	5.0	7.0	5.0	ND	5.0	25	5.0
Zinc	ICAP	ND	5.0	ND	5.0	ND	5.0	27	5.0	25	5.0	ND	5.0

Table 1.23 (Cont) Results of the Metals Analysis for the Water Samples WA # 0-098 Pier Drum

Client ID Location		B00641 Filtered		A00635 Unfiltere	ed	B00635 Filtered		A00639 Unfiltere	d	B00639 Filtered		A00640 Unfiltere	ed .
Parameter	Analysis Method	Conc ug/l	MDL ug/l	Conc ug/l		Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l
Aluminum	ICAP	3400	40	1400	40	1100	40	2000	40	1400	40	ND	40
Antimony	AA-Fur	ND	5.6	ND	5.6	ND	5.6	ND	5.6	ND	5.6	ND	5.6
Arsenic	AA-Fur	10	2.2	3.0	2.2	2.8	2.2	10	2.2	9.6	2.2	ND	2.2
Barium	ICAP	4.0	4.0	18	4.0	17	4.0	9.0	4.0	7.0	4.0	ND	4.0
Beryllium	ICAP	ND	2.0	ND	2.0	ND	2.0	ND	2.0	ND	2.0	ND	2.0
Cadmium	AA-Fur	ND	0.22	ND	0.22	ND	0.22	ND	0.22	ND	0.22	ND	0.22
Calcium	ICAP	1700	100	4100	100	4100	100	12000	100	9800	100	ND	100
Chromium	AA-Fur	ND	2.8	ND	2.8	ND	2.8	3.6	2.8	2.9	2.8	ND	2.8
Cobalt	ICAP	ND	6.0	ND	6.0	ND	6.0	ND	6.0	ND	6.0	ND	6.0
Copper	ICAP	ND	5.0	ND	5.0	ND	5.0	ND	5.0	ND	5.0	ND	5.0
Iron	ICAP	1700	10	9200	10	9200	10	1600	10	1100	10	ND	10
Lead	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2
Magnesium	ICAP	870	500	4300	500	4200	500	4600	500	3600	500	ND	500
Manganese	ICAP	ND	2.0	3.0	2.0	3.0	2.0	26	2.0	21	2.0	ND	2.0
Mercury	Cold Vapor	ND	0.20	ND	0.20	ND	0.20	ND	0.20	ND	0.20	/ND	0.20
Nickel	ICAP	ND	10	ND	10	ND	10	ND	10	ND	10	ND	10
Potassium	ICAP	ND	2000	ND	2000	ND	2000	2300	2000	ND	2000	ND	2000
Selenium	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2
Silver	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2
Sodium	ICAP	3000	500	30000	500	30000	500	30000	500	23000	500	ND	500
Thallium	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2
Vanadium	ICAP	9.0	5.0	ND	5.0	ND	5.0	8.0	5.0	6.0	5.0	ND	5.0
Zinc	ICAP	ND	5.0	ND	5.0	ND	5.0	56	5.0	45	5.0	ND	5.0

Table 1.23 (Cont) Results of the Metals Analysis for the Water Samples WA # 0-098 Pier Drum

Client ID Location		B00640 Filtered		A00634 Unfilters	d	B00634 Filtered		A00637 Unfilters	d	B00637 Filtered	
Parameter	Analysis Method	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/l
Aluminum	ICAP	ND	40	7400	40	3500	40	10000	40	4100	40
Antimony	AA~Fur	ND	5.6	ND	5.6	ND	5.6	ND	5.6	ND	5.6
Arsenic	AA-Fur	ND	2.2	10	2.2	10	2.2	19	2.2	15	2.2
Barium	ICAP	ND	4.0	30	4.0	10	4.0	39	4.0	9.0	4.0
Beryllium	ICAP	ND	2.0	ND	2.0	ND	2.0	ND	2.0	ND	2.0
Cadmium	AA-Fur	ND	0.22	ND	0.22	ND	0.22	ND	0.22	ND	0.22
Calcium	ICAP	ND	100	27000	100	28000	100	5800	100	6000	100
Chromium	AA~Fur	ND	2.8	7.6	2.8	3.8	2.8	13	2.8	3.5	2.8
Cobatt	ICAP	ND	6.0	ND	6.0	ND	6.0	ND	6.0	ND	6.0
Copper	ICAP	ND	5.0	ND	5.0	ND	5.0	ND	5.0	ND	5.0
Iron	ICAP	ND	10	6600	10	5800	10	4500	10	3600	10
Lead	AA-Fur	ND	2.2	ND	2.2	ND	2.2	3.2	2.2	ND	2.2
Magnesium	ICAP	ND	500	13000	500	13000	500	5800	500	5600	500
Manganese	ICAP	ND	2.0	6.0	2.0	3.0	2.0	8.0	2.0	5.0	2.0
Mercury	Cold Vapor	ND	0.20	ND	0.20	ND	0.20	ND	0.20	ND	0.20
Nickel	ICAP	ND	10	ND	10	ND	10	ND	10	ΝĐ	10
Potassium	ICAP	ND	2000	2100	2000	ND	2000	ND	2000	ND	2000
Selenium	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2
Silver	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2
Sodium	ICAP	ND	500	75000	500	76000	500	40000	500	40000	500
Thallium	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	ND	2.2
Vanadium	ICAP	ND	5.0	12	5.0	7.0	5.0	11	5.0	ND	5.0
Zinc	ICAP	ND	5.0	ND	5.0	5.0	5.0	ND	5.0	ND	5.0

1

Table 1.24 Results of the Metals Analysis for the Soil Samples WA # 0-098 Pier Drum
Based on Dry Weight

2 2 3362

Client ID		Method	Blank	A00642		A00643		A00644	
Location		-		-		-		-	
Parameter	Analysis Method	Conc mg/kg	MDL mg/kg	Conc mg/kg	MDL mg/kg	Conc mg/kg	MDL mg/kg	Conc mg/kg	MDL mg/kg
Aluminum	ICAP	ND	10	660	9.4	470	8.9	330	9.3
Antimony	AA-Fur	ND	1.0	ND	0.84	ND	0.89	ND	0.83
Arsenic	AA-Fur	ND	0.50	ND	0.42	ND	0.45	ND	0.42
Barium	ICAP	ND	4.0	4.7	3.8	3.6	3.6	ND	3.7
Beryllium	ICAP	ND	0.2	ND	0.2	ND	0.2	ND	0.2
Cadmium	ICAP	ND	0.3	ND	0.3	ND	0.3	ND	0.3
Calcium	ICAP	ND	50	290	47	94	45	61	47
Chromium	ICAP	ND	0.8	1.4	0.8	ND	0.7	ND	0.7
Cobalt	ICAP	ND	2.0	ND	1.9	ND	1.8	ND	1.9
Copper	ICAP	ND	0.6	ND	0.6	ND	0.5	ND	0.6
Iron	ICAP	ND	9.0	200	8.4	150	8.0	43	8.4
Lead	ICAP	ND	4.0	ND	3.8	ND	3.6	ND	3.7
Magnesium	ICAP	ND	50	ND	47	ND	45	ND	47
Manganese	ICAP	ND	1.4	ND	1.3	ND	1.3	ND	1.3
Mercury	Cold Vapor	ND	0.04	ND	0.03	ND	0.03	ND	0.04
Nickel	ICAP	ND	2.0	ND	1.9	ND	1.8	ND	1.9
Potassium	ICAP	ND	200	ND	190	ND	180	ND	190
Selenium	AA-Fur	ND	0.50	ND	0.42	ND	0.45	ND	0.42
Silver	ICAP	ND	0.5	ND	0.5	ND	0.4	ND	0.5
Sodium	ICAP	ND	50	69	47	63	45	93	47
Thallium	AA-Fur	ND	0.50	ND	0.42	ND	0.45	ND	0.42
Vanadium	ICAP	ND	2.0	ND	1.9	ND	1.8	ND	1.9
Zinc	ICAP	ND	2.0	5.4	1.9	ND	1.8	3.4	1.9

Table 1.25 Results of the TCLP Metals Analysis WA # 0-098 Pier Drum

Client ID Location		Method Lab	Blank	A00525 A00642		A00526 A00643		A,D00527 A00644		
Parameter	Analysis Method	Conc ug/i	MDL ug/l	Conc ug/l	MDL ug/l	Conc ug/l	MDL ug/i	Conc ug/l	MDL ug/i	
Arsenic	AA-Fur	ND	2.2	6.3	2.2	ND	2.2	ND	2.2	
Barium	ICAP	ND	4.0	930	4.0	28	4.0	11	4.0	
Cadmium	ICAP	ND	3.0	ND	3.0	ND	3.0	ND	3.0	
Chromium	ICAP	ND	5.0	8.0	5.0	ND	5.0	ND	5.0	
Lead	AA-Fur	ND	2.2	3.2	2.2	ND	2.2	ND	2.2	
Mercury	Cold Vapor	ND	0.20	ND	0.20	ND	0.20	ND	0.20	
Selenium	AA-Fur	ND	2.2	ND	2.2	ND	2.2	ND	2.2	
Silver	ICAP	ND	10	ND	10	ND	10	ND	10	

Table 1.26 Results of the Metals Analysis for the Drum Liquid Samples WA # 0-098 Pier Drum

Client ID Location		Method Lab	Blank	00912 Reac#9		00913 Reac#2		00914 Reac#4	(°) •
Parameter	Analysis	Conc	MDL	Conc	MDL	Conc	MDL	Conc	MDL
	Method	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l
Aluminum Antimony Arsenic Barium Beryllium Cadmium Calcium Chromium Cobalt Copper Iron Lead Magnesium Manganese Nickel Potassium Selenium Silver	ICAP AA-Fur ICAP ICAP		50 5.0 5.0 2.0 2.0 100 5.0 5.0 5.0 25 5.0 200 200 5.0	690 ND ND 29 ND 7.0 410* 80 7300 520 18* 61 8000 3000 66 12000 ND ND	100 10 10 4.0 4.0 0.20 * 10 10 0.05 * 10 1000 4.0 20 4000 10	ND ND ND 34 400* 120 40000 150 360* 1200 8000 8000 8500 850 35000 ND ND	100 10 10 4.0 4.0 0.20 * 10 10 0.05 * 1000 4.0 20 4000 10	47 900 660 310* 38 10000 1500 30 12000 ND ND	100 10 10 10 4.0 4.0 0.20 * 10 10 0.05 * 10 1000 4.0 20 4000 10 10
Sodium	ICAP	ND	500	21000	1000	71000	1000	29000	1000
Thallium	AA-Fur	ND	5.0	ND	10	ND	10	ND	10
Vanadium	ICAP	ND	5.0	ND	10	ND	10	32	10
Zinc	ICAP	ND	5.0	1900	10	3600	10	330	10

J364

^{*} denotes that the units for this value is milligram per liter Mercury analysis were not performed due to insufficient sample size

Table 1.27 Results of the Metals Analysis in TCLP Extracts from the Drum Solids WA # 0-098 Pier Drum

Client ID Location	Method Bis	enk	00911 Composit 1, 8, 9, 11		00915 REAC #5		00916 Composit 3, 7, 10, 13, 14		
Parameter	Conc (ug/l)	MDL (ug/l)	Conc (ug/l)	MDL (ug/l)	Conc (ug/l)	MDL (ug/l)	Conc (ug/l)	MDL (ug/l)	
Aluminum	ND	3000	ND	3000	ND	3000	ND	3000	
Antimony	ND	10.0	ND	10.0	ND	10.0	ND	10.0	
Arsenic	ND	8.0	ND	8.0	ND	8.0	ND	8.0	
Bartum	ND	1000	ND	1000	ND	1000	ND	1000	
Beryllium	ND	1.0	ND	1.0	ND	1.0	ND	1.0	
Cadmium	ND	4.0	167	80	ND	4.0	ND	4.0	
Calcium	ND	250	435000	25000	2800	250	1290	250	
Chromium	ND	100	ND	100.0	ND	100.0	ND	100.0	
Cobalt	ND	250	470	250	ND	250	ND	250	
Copper	ND	50.0	ND	50.0	ND	50.0	50	50.0	
Iron	ND	200	2190	200	ND	200	760	200	
Lead	ND	5.0	102	20	7.9	5.0	6.1	5.0	
Magnesium	ND	50.0	4300	500	300	50.0	170	50.0	
Manganese	ND	100	350	100	ND	100	ND	100	
Mercury	ND	0.5	ND	0.5	ND	0.5	ND	0.5	
Nickel	ND	100	ND	100	ND	100	ND	100	
Potassium	ND	100	4200	1000	1580	100	570	100	
Selenium	ND	10.0	ND	10.0	ND	10.0	ND	10.0	
Silver	ND	50.0	ND	50.0	ND	50.0	ND	50.0	
Sodium	ND	50.0	3100000	500000	2300000	500000	1800000	500000	
Thallium	ND	10.0	ND	10.0	ND	10.0	ND	/ 10.0	
Vanadium	ND	3000	ND	3000	ND	3000	ND	3000	
Zinc	ND	50.0	440	50.0	210	50.0	130	50.0	

Table 1.28 Results of the Wet Chemistry Analysis WA # 0~098 Pier Drum

Sample ID	Location	Analysis	Units	Result	MDL
00912	REAC #9	вти	BTU/lb	3379	50
00912	REAC #9	Ash	Percent	0.2	0.1
00912	REAC #9	CI	Percent	0.33	0.001
00912	REAC #9	S	Percent	ND	0.1
00912	REAC #9	Water	Percent	53	1
00913	REAC #2	BTU	BTU/lb	5470	50
00913	REAC #2	Ash	Percent	0.26	0.1
00913	REAC #2	a	Percent	0.34	0.001
00913	REAC #2	S	Percent	ND	0.1
00913	REAC #2	Water	Percent	47	1
00914	REAC #4	BTU	BTU/Ib	ND	50
00914	REAC #4	Ash	Percent	ND	0.1
00914	REAC #4	CI	Percent	0.44	0.001
00914	REAC #4	S	Percent	0.10	0.1
00914	REAC #4	Water	Percent	89.6	1

2 2 3367

QA/QC FOR VOA

Before extraction, each sample was spiked with a three component mixture of CLP surrogate standards consisting of toluene-d_a, 4-bromofluorobenzene and 1,2-dichloroethane-d₄. The surrogate percent recoveries for the water samples, listed in Table 2.1, ranged from 90 to 102. All forty-five values were within the acceptable QC limits.

The internal standard areas (for bromochloromethane, 1,4-difluorobenzene, and chlorobenzene-d₅) for the water samples are also listed in Table 2.1. All forty-five areas are within acceptable QC limits.

Sample 00635 was chosen for the matrix spike/matrix spike duplicate (MS/MSD) analyses for the water samples. The percent recoveries, ranging from 94 to 110, are listed in Table 2.2. All ten values were within the acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.2, ranged from 0 (zero) to five, and all five were within the acceptable QC limits.

The surrogate percent recoveries for the soil samples, listed in Table 2.3, ranged from 86 to 105. All twenty-four values were within acceptable QC limits.

The internal standard areas for the soil samples are also listed in Table 2.3. All twenty-four areas are within the acceptable QC limits.

Sample 00642 was chosen for the MS/MSD analyses for the soil samples. The percent recoveries, ranging from 99 to 106, are listed in Table 2.4. All ten values were within the acceptable QC limits. The RPDs, also listed in Table 2.4, ranged from 1 to 3, and all five were within the acceptable QC limits.

The surrogate percent recoveries for the drum liquid samples, listed in Table 2.5, ranged from 93 to 105. All thirty-six values were within acceptable QC limits.

The internal standard areas for the drum liquid samples are also listed in Table 2.5. All thirty-six areas are within the acceptable QC limits.

Sample 00914 was chosen for the MS/MSD analyses for the drum liquid samples. The percent recoveries, ranging from 89 to 106, are listed in Table 2.6. All ten values were within the acceptable QC limits. The RPDs, also listed in Table 2.6, ranged from 0 (zero) to 8, and all five were within the acceptable QC limits.

The initial calibrations are listed in Table 2.7.

The continuing calibrations are listed in Table 2.8.

The samples and calibrations are correlated in Table 2.9.

The surrogate percent recoveries for the TCLP samples, listed in Table 2.10, ranged from 98 to 115. Nineteen out of twenty-one values were within the acceptable QC limits.

The internal standard areas for the TCLP samples are listed in Table 2.11. All eighteen areas are within the acceptable QC limits.

Non-REAC sample 95-01-227-1 was chosen for the MS/MSD analyses for the TCLP samples. The percent recoveries, ranging from 64 to 86, are listed in Table 2.12. All ten values were within the acceptable QC limits. The RPDs, also listed in Table 2.12, ranged from 0 (zero) to 6, and all five were within the acceptable QC limits.

The surrogate percent recoveries for the TCLP extracts of the drum solid samples, listed in Table 2.13, ranged from 79 to 111. All twenty-one values were within the acceptable QC limits.

The internal standard areas for the TCLP extracts of the drum solid samples are listed in Table 2.14. All twenty-four values are within the acceptable QC limits.

The blank spike percent recoveries, listed in Table 2.15, ranged from 90 to 140. Four out of five values are within the acceptable QC limits.

Sample 00911 was chosen for the MS/MSD analyses for the TCLP extracts from the drum solid samples. The percent recoveries, ranging from 97 to 115, are listed in Table 2.16. All ten values were within the acceptable QC limits. The RPDs, also listed in Table 2.16, ranged from 1 to 4, and all five values were within the acceptable QC limits.

Table 2.1 Results of the Internal Standard Areas and Surrogate Recoveries for the Water Samples WA # 0-098 Pier Drum

Sample # File area area area 270346 NA NA LAB BLANK-01/13/95 >V2426 75692 314945 273790 94 97 00645 >V2427 76089 320731 284218 96 99 00638 >V2430 73332 306255 275550 98 97 00635MSD >V2436 74683 305104 261234 100 101 00635MSD >V2437 69724 284515 244748 102 101 CAL CHECK 50 PPB VOC >V2440 71286 311633 263301 NA NA LAB BLANK-01/14/95 >V2441 71588 316662 268936 90 99 00640 >V2443 71161 302822 264733 99 97
CAL CHECK 50 PPB VOC >V2425 76546 321010 270346 NA NA LAB BLANK-01/13/95 >V2426 75692 314945 273790 94 97 00645 >V2427 76089 320731 284218 96 99 00638 >V2430 73332 306255 275550 98 97 000635MS >V2436 74683 305104 261234 100 101 00635MSD >V2437 69724 284515 244748 102 101 CAL CHECK 50 PPB VOC >V2440 71286 311633 263301 NA NA LAB BLANK-01/14/95 >V2441 71588 316662 268936 90 99
LAB BLANK-01/13/95 >V2426 75692 314945 273790 94 97 00645 >V2427 76089 320731 284218 96 99 00638 >V2430 73332 306255 275550 98 97 00635MS >V2436 74683 305104 261234 100 101 00635MSD >V2437 69724 284515 244748 102 101 CAL CHECK 50 PPB VOC >V2440 71286 311633 263301 NA NA LAB BLANK-01/14/95 >V2441 71588 316662 268936 90 99
00645
00638
CAL CHECK 50 PPB VOC >V2440 71286 311633 263301 NA NA AB BLANK-01/14/95 >V2441 71588 316662 268936 90 99
AB BLANK-01/14/95 >V2441 71588 316662 268936 90 99
00640 >v2443 71161 302822 264733 99 97
CAL CHECK 50 PPB VOC >V2467 69611 297158 256657 NA NA
AB BLANK-01/17/95 >V2472 69211 297347 253604 96 97
00636 >v2473 70573 300017 255424 95 99
00641 >V2474 66183 276961 240465 96 97
00635 >v2475 70963 303151 259143 96 97
00634 >v2477 69445 295544 248081 95 98
00637 >v2478 65921 286093 242340 96 98
00637 >V2478 65921 286093 242340 96 98

SURROGATE LIMITS

WATER

^{\$1 (}DIC) = 1,2-Dichloroethane-d4 (76-114) \$2 (TOL) = Toluene-d8 (88-110) \$3 (BRO) = Bromofluorobenzene (86-115)

2 2 33/4

Table 2.2 Results of the MS/MSD Analysis for the Water Samples WA # 0-098 Pier Drum

Matrix Spike - EPA Sample No.: 00635

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS % REC #	QC LIMITS REC.
=======================================	========	2222222222	**=**=**=	=====	======
1,1-Dichloroethene	50.00	0.00	47.23	94	61-145
Trichloroethene	50.00	0.00	53.89	108	71-120
Benzene	50.00	0.00	51.18	102	76-127
Toluene	50.00	0.00	49.45	99	76-125
Chlorobenzene	50.00	0.00	49.49	99	75 - 130

	SP1KE ADDED	MSD CONCENTRATION	MSD %	x	QC L	IMITS
COMPOUND	(ug/L)	(ug/L)	REC #	RPD #	RPD	REC.
	========			222322	======	=====
1,1-Dichloroethene	50.00	49.55	99	5	14	61-145
Trichloroethene	50.00	55.21	110	2	14	71-120
Benzene	50.00	53.41	107	5	11	76-127
Toluene	50.00	51.35	103	4	13	76-125
Chlorobenzene	50.00	49.61	99	0	13	75-130
				i	İ	İ

Table 2.3 Results of the Internal Standard Areas and Surrogate Recoveries for the Soil Samples WA # 0-098 Pier Drum

			Inter	nal Standard	ls	Surrogates			
CAL CHECK 5	Sample # 50 PPB VOC	Data File >V2440	1 area 71286	2 area 311633	3 area 263301	DIC % NA	TOL % NA	BRO % NA	
SAND BLANK-C	1/14/95	>V2442	58256	239724	216206	9 5	98	99	
00	0643	>V2451	62773	273702	209206	99	98	97	
00	0644	>V2452	62044	269343	222593	101	98	97	
00	0642MS	>V2453	60492	257058	203472	103	102	96	
00	642MSD	>V2454	60973	259847	206304	104	105	98	

CAL CHECK 50 PPB VOC	>V2459	72261	306490	262689	NA	NA	NA
SAND BLANK-01/17/95	>V2460	57591	245235	206015	95	99	95
00642	>V2461	61869	268131	221075	95	103	86
(1:5 DIL) 00642	>V2462	67440	295631	250991	98	99	95

SURROGATE LIMITS

SOIL

S1 (D1C) = 1,2-Dichloroethane-d4 (70-121) S2 (TOL) = Toluene-d8 (81-117) S3 (BRO) = Bromofluorobenzene (74-121)

Table 2.4 Results of the MS/MSD Analysis for the Soil Samples WA # 0-098 Pier Drum

Matrix Spike - EPA Sample No.: 00642

Level:(low/med) LOW

COMPOUND	SPIKE	SAMPLE	MS	MS	QC
	ADDED	CONCENTRATION	CONCENTRATION	%	LIMITS
	(ug/Kg)	(ug/Kg)	(ug/Kg)	REC #	REC.
1,1-Dichloroethene	63.29 63.29 63.29 63.29 63.29	0.00 0.00 11.90	66.34 64.91 76.92	105 103	59-172 62-137 66-142 59-139 60-133

	SPIKE ADDED	MSD CONCENTRATION	MSD %	x	QC L	IMITS
COMPOUND	(ug/Kg)	(ug/Kg) ·	REC #	RPD #	RPD	REC.
=======================================	=======	=========	=====	======	=====	=====
1,1-Dichloroethene	63.29	63.75	101	1	22	59-172
Trichloroethene	63.29	67.16	106	1	24	62-137
Benzene	63.29	64.01	101	2	21	66-142
Toluene	63.29	75.42	100	3	21	59-139
Chlorobenzene	63.29	63.24	100	1	21	60-133
				İ	i	į

2 2 33/3

Table 2.5 Results of the Internal Standard Areas and Surrogate Recoveries for the Drum Liquid Samples WA # 0.098 Pier Drum

			Inter	nal Standard	is	Sur	rogates	;
CAL CHECK	Sample #	Data File >V2618	1 area 51926	2 area 21 7 627	3 area 180758	DIC % NA	TOL % NA	BRO % NA
LAB BLANK-	01/30/95	>V2619	49694	215774	182052	101	98	98
	00914	>V2620	51010	225513	188614	100	98	98
	00912	>V2621	48851	220310	178397	103	99	98
	00913	>V2622	48472	212162	180851	101	96	96
	00914MS	>V2624	49468	219840	177481	105	101	101
	00914MSD	>V2625	50118	224368	178615	103	102	100
CAL CHECK	50 PPB VOC	>v2628	52187	222379	187140	NA	NA	NA
LAB BLANK-	01/31/95	>V2629	52983	215667	194137	96	98	97
500000DL	00912	>V2631	45750	216746	179827	97	100	95
500000DL	00913	>v2632	47964	218820	182803	98	99	97
20DL	00914	>v2633	49769	219375	183169	9 5	99	96
1000000DL	00912	>v26 3 4	42135	223211	163060	97	101	93
2500000DL	00913	>v26 3 5	47524	211477	178429	97	100	96

SURROGATE LIMITS

WATER

S1 (DIC) = 1,2-Dichloroethane-d4 (76-114) S2 (TOL) = Toluene-d8 (88-110) S3 (BRO) = Bromofluorobenzene (86-115)

2 : 33/

Table 2.6 Results of the MS/MSD Analysis for the Drum liquid Samples WA # 0-098 Pier Drum

Matrix Spike - EPA Sample No.: 00914

	SPIKE	SAMPLE	MS	MS	90
,	ADDED	CONCENTRATION	CONCENTRATION	*	LIMITS
COMPOUND	(ug/L)	(ug/L)	(ug/L)	REC #	REC.
	=======	=======================================		=====	=====
1,1-Dichloroethene	50.00	0.00	52.49	105	61-145
Trichloroethene	50.00	0.00	47.87	96	71-120
Benzene	50.00	0.00	52.07	104	76-127
Toluene	50.00	9.58	60.66	102	76-125
Chlorobenzene	50.00	0.00	50.42	101	75-130

	SPIKE ADDED	MSD CONCENTRATION	MSD %	x	QC L	IMITS
COMPOUND	(ug/L)	(ug/L)	REC #	RPD #	RPD	REC.
=======================================	*******	=======================================	=====	2==2==	======	
1,1-Dichloroethene	50.00	52.65	105	0	14	61-145
Trichloroethene	50.00	44.67	89	8	14	71-120
Benzene	50.00	53.11	106	2	11	76-127
Toluene	50.00	61.56	104	2	13	76-125
Chlorobenzene	50.00	51.29	103	2	13	75-130
						!

Table 2.7 Results of the Initial Calibration WA # 0-098 Pier Drum

Instrument ID: GCMSD-2(3034A12982)

Calibration Date: 01/18/95

Minimum RF for SPCC is 0.30

Maximum % RSD for CCC is 30.0%

Laboratory II	0: >V2465 RF 5.00	>V2466 RF 20.00	>V2467 RF 50.00	>V2468 RF 100.00	>V2469 RF 150.00	>V2470 RF 200.00	RRT	RF	% RSD	CCC SPCC
Dichlorodifluoromethane	2.80941	2.50632	2.34395	2.33993	2.18946	2.27145	.233	2.41009	9.202	
Chloromethane				1.01467		1.00117		1.04526	8.101	**
Vinyl Chloride	1.19594	1.08058	1.01077	1.00967	.95248		.273	1.03762	8.558	*
Bromomethane		1.19411			.83452			1.07394	24.028	
Chloroethane	.78867	.73370	.68360	.65068	.41482	.42918	.348	.61678	25.617	•
Trichlorofluoromethane	4.37044	3.86030	3.68695	3.70158			.401	3.77576	8,501	
Acetone	.57416	.48424	.46921	.43028	.43098	. 42623	.496	.46918	12.081	
1,1-Dichloroethene	1.57402	1.38744	1.33171	1.30759	1.21965	1.26641	.515	1.34780	9.247	*
Carbon Disulfide	3.34714	3.13399	3.12541	3.19138	3.03479	3.15376		3.16441	3,268	
Methylene Chloride				1.33106	-			1.39021	11.135	
rans-1,2-Dichloroethene	1.70857	1.51001	1.46869	1.45634	1.36206	1.37998	.701	1.48094	8.420	
lethyl-tertiary-butylether				2.61744	_			2.75552	6.992	
,1-Dichloroethane	3.00942	2.78345	2.68241	2.72711	2.53989	2.61298	.800	2.72588	5.982	**
2-Butanone	.06878	.10413	.10524	.09518	.11644	.11347	.918	.10054	17.182	
.2-Dichloropropane	2.28892	2.08137	2.05150	2.09128	2.02741	2.10641	.941	2.10782	4,421	
is-1,2-Dichloroethene	1.73321	1.61938	1.51411	1.49753	1.41357	1.42604	.941	1.53397	7.977	
hloroform	3.63603	3.34706	3.22257	3.24602	3.09260	3.10819	.982	3.27541	6,109	*
.1-Dichloropropene	2.25944	2.14454	2.09126	2.04245	1.96480	1.97036	1.114	2.07881	5.408	
,2-Dichloroethane				1.97685			1,154	2.05552	7.420	
.2-Dichloroethane-d4 (SURR)	1.72176	1.74948	1.79147	1.78923	1.81991	1.84581	1.133	1.78628	2.532	(Conc=50.0,5
,1,1-Trichloroethane	.66742	.70216	.69706	.70441	.70565	.72301	.860	.69995	2.600	•
arbon Tetrachloride	.34904	.44551	.50682	.55326	.57513	.59830	.904	.50468	18,549	
enzene	.87507	.86867	.84555	.82848	.81981	.82275	.926	.84339	2,832	
richloroethene	.49510		.43654	.41657	.43129	.43271	1.047	.44451	6.223	
,2-Dichloropropane	.42611	.35643	.38555	.39837	.33326	.37312	1.073	.37881	8.573	*
ibromomethane	. 44068	.43096	.41777	.40334	.39685	.39163	1.106	.41354	4.739	
romodichloromethane	.61514	.65462	.69746	.69728	.69515	.70348	1.111	.67719	5.195	
is-1,3-Dichtoropropene	.46703		.48387	.47599	.47396	.47046	1.229	.47831	2.392	
rans-1,3-Dichloropropene	.35836	.41618	.42532	.41177	.40916	.40699	1.328	.40463	5.827	
1,1,2-Trichloroethane	. 29021	.31408	.31045	. 29529	.30172	.29266	1.349	.30073	3.255	

RF - Response Factor (Subscript is amount in ppb)

RRT - Average Relative Retention Time (RT Std/RT Istd)

RF - Average Response Factor

%RSD - Percent Relative Standard Deviation CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

2 1 3376

Table 2.7 (Cont) Results of the Initial Calibration WA # 0-098 Pier Drum

Instrument ID: GCMSD-2(3034A12982) Calibration Date: 01/18/95

Minimum RF for SPCC is 0.30

Maximum % RSD for CCC is 30.0%

Laboratory ID:	RF	>V2466 RF	>V2467 RF	>V2468 RF	>V2469 RF	>V2470 RF		RF	* 000	000 0000
Compound	5.00	20.00	50.00	100.00	150.00	200.00	RRT	Kr	% RSD	CCC SPCC
1,3-Dichloropropane	.50841	.53170	.53208	.50405	.49318	.48497	1.398	.50907	3.830	
Dibromochloromethane	.57879	.64674	.69595	.66219	.67848	.66363	1.428	.65430	6.200	
1,2-Dibromoethane	.47894	.52921	.54458	.51182	.50526	.48724	1.465	.50951	4.865	
Bromoform	.38227	.48071	.53641	.51264	.55750	.53781	1.721	.50122	12.758	**
4-Methyl-2-Pentanone	.10495	.10996	.11177	.10424	.11298	.11538	.783	.10988	4.062	•
Toluene-d8 (SURR)		1.19282				1.24098		1.20579		(Conc≈50.0,50.0,5
Toluene	.83660	.79838	.76866	.76718	.76298	.79133	.830	.78752	3.552	*
2-Hexanone	.12050	.13015	.12865	.11839	.12389	.12467	.890	.12437	3.644	
Tetrachloroethene	.78176	.74764	.73977	.72982	.71381	.73934	.910	.74202	3.054	
Chlorobenzene	1.12423	1.06312	1.06099	1.02727	1.02361	1.05245	1.005	1.05861	3.426	**
1,1,1,2-Tetrachloroethane	.55853	.56901	.59500	.58501	.58870	.61335	1.014	.58493	3.309	
Ethylbenzene	1.80006	1.75738	1.73153	1.71226	1.71240	1.76644	1.022	1.74668	1.974	*
p & m-Xylene	1.58425	1.46123	1.47086	1.42952	1.31695	1.22154	1.034	1.41406	9.009	(Conc=10.0,40.0,1
o-Xylene	1.56453	1.43732	1.43031	1.38335	1.36407	1.41104	1.084	1.43177	4.940	
Styrene	.66587	.65324	.64826	.63423	.62457	.64754	1.087	.64562	2.246	
Isopropylbenzene	1.97084	1.90899	1.89950	1.85722	1.83039	1.90184	1.138	1.89480	2.544	
1,1,2,2-Tetrachloroethane	.54403	.60821	.59361	.56082	.57918	.57530	1.150	.57686	3.958	**
p-Bromofluorobenzene (SURR)	.66259	.67274	.69490	.67076	.67236	.67872	1.157	.67535	1.613	(Conc≈50.0,50.0,5
1,2,3-Trichloropropane	. 16595	.17142	.17100	. 16278	. 16796	. 16700	1.166	.16768	1.933	
Bromobenzene	.68284	.63855	.64685	.61595	.61440	.62785	1.173	.63774	3.989	
n-Propylbenzene	.49170	.46927	.45507	.44126	. 42975	.44441	1.190	.45524	4.900	
2-Chlorotoluene	. 48663	.42607	.40294	.39938	.38532	.40550	1.199	.41764	8.681	
4-Chlorotoluene	.48214	.45660	.48579	.44150	.43873	.44340	1.208	.45803	4.594	
1,3,5-Trimethylbenzene				1.56987				1.62309	4.350	
tert-Butylbenzene				1.55561				1.61365	4.473	
1,2,4-Trimethylbenzene				1.52618				1.56897	4.252	
sec-Butylbenzene				2.16661				2.21808	3.628	
1,3-Dichlorobenzene		1.04545		.99940		1.01295		1.03711	5.365	
p-Isopropyltoluene				1.79425				1.80992	3.385	
1,4-Dichlorobenzene				1.08151				1.12622	5.326	
1,2-Dichlorobenzene	1.07153		.99370	.93613	.93850		1.357		5.232	
n-Butyl benzene				1.80177				1.84422	4.377	
1,2-Dibromo-3-Chloropropane	.10564	.13720		. 13244	. 14558	.14385	1.458		11.024	
1,2,4-Trichlorobenzene	.95844	.91394	.92259		.87825	.88889	1.583	.90669	3.463	
Naphthalene Hexachlorobutadiene	.95034	.85899	.86401	1.09341 .82592	1.16488	1.16419	1.616	1.12938	2.685 5.375	
1,2,3-Trichlorobenzene	.89627			.76448	.78040		1.640		6.087	
1,2,3 II ICIICOI ODENZENE	.07021	.03547		. / 0-40	. / 0040	.,,,,,	1.040	.00740	0.007	

RF - Response Factor (Subscript is amount in ppb)

RRT - Average Relative Retention Time (RT Std/RT Istd)

RF - Average Response factor

XRSD - Percent Relative Standard Deviation CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

098\DEL\AR\9503\QAQCVOA

Table 2.7 (Cont) Results of the Initial Calibration WA # 0-098 Pier Drum

Instrument ID: GCMSD-2(3034A12982) Calibration Date: 11/22/94

Minimum RF for SPCC is 0.30

Maximum % RSD for CCC is 30.0%

Laboratory ID	: >V2218 RF 5.00	>V2219 RF 20.00	>V2220 RF 50.00	>V2221 RF 100.00	>V2222 RF 150.00	>V2223 RF 200.00	RRT	RF	% RSD	ccc cpcc
Compound	3.00	20.00	50.00	100.00	130.00	200.00	KKI	Kr	A KSU	CCC SPCC
Dichlorodifluoromethane	3 35896	2.81799	2.76825	2,61334	2.50108	2.47974	. 232	2.75656	11,802	
Chloromethane				1.06640				1.10099	9,408	**
Vinyl Chloride				1.02462		.96592		1.06518	10.066	•
Bromomethane		1.13568			.81755	.82885	_	1.00911	22.225	
Chloroethane	.78744		.69770				.346		25.910	
Trichlorofluoromethane				3.56132				3.68991	10.269	•
Acetone	.38773					.23102	.493		19.013	
1,1-Dichloroethene				1.27812				1.31810	11.099	*
Carbon Disulfide				3.10763				3.09811	5.084	
Methylene Chloride				1.30529				1.38705	13.679	
trans-1,2-Dichloroethene				1.38464				1.42359	9.076	
Methyl-tertiary-butylether				2.74496				2.76397	13.194	
1.1-Dichloroethane				2.67014				2.74346	8.653	**
2-Butanone	.07208	.08411	.08000			.09193	.919	.08201	11.024	
2,2-Dichloropropane				2.02129				2.03221	5.480	
cis-1.2-Dichloroethene				1.46920				1.50057	9.205	
Chloroform				3.15613				3.27645	10.321	*
1,1-Dichloropropene				1.95511				2.03304	8.425	
1.2-Dichloroethane				2.04983				2.12352	10.741	
1,2-Dichloroethane-d4 (SURR)				1.87296				1.84429		(Conc=50.0,50.0,
1,1,1-Trichloroethane	. 78193	.68049	.67570		.69106	.70636	.861	.70440	5.597	,
Carbon Tetrachloride	.42121	.43781	.49230		.54085	.55433	.904	.49450	11.074	
Benzene	1.02658	.85071	.83508	.82616	.81930	.82015	.927	.86300	9.384	
Trichloroethene	.53170	.46562	.41950	.39292	.42206	.45497	1.048	.44779	10.887	
1,2-Dichloropropane	.48560	.41776	.39434	.40416	.40117	.40755	1.075	.41843	8.078	*
Dibromomethane	.52348	.42565	.41265	.39983	.38500	.38811	1.108	.42245	12.257	
Bromodichloromethane	.72935	.66218	.69006	.70695	.69877	.70450	1.112	.69864	3.168	
cis-1,3-Dichloropropene	.56723	.47345	.46314	.47820	.45994	.45670	1.232	.48311	8.697	
trans-1,3-Dichloropropene	.46088	.42743	.41911	.41528	.40101	.39057	1.328	.41905	5.817	
1,1,2-Trichloroethane	.36864	.32150	.31471	.31380	.30360	.29777	1.350	.32000	7.903	

RF - Response Factor (Subscript is amount in ppb)

RRT - Average Relative Retention Time (RT Std/RT Istd)

RF - Average Response Factor

%RSD - Percent Relative Standard Deviation CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

Table 2.7 (Cont) Results of the Initial Calibration WA # 0-098 Pier Drum

Instrument ID: GCMSD-2(3034A12982)

Calibration Date: 11/22/94

Minimum RF for SPCC is 0.30 Maximum % RSD for CCC is 30.0%

Laboratory [Compound	D: >V2218 RF 5.00	>V2219 RF 20.00	>V2220 RF 50.00	>V2221 RF 100.00	>V2222 RF 150.00	>V2223 RF 200.00	RRT	RF	% RSD	CCC SPCC
1.7.Dishlaramana	.64190	.54413	.53676	.52963	.49669	.48207	1.398	.53853	10.417	•••
1,3-Dichloropropane Dibromochloromethane	.70026		.64435	.63321	.63975	.63574	1.428	.64683	4.141	
1,2-Dibromoethane	.61299		.53295	.53397	.49539	.47095	1.465	.52909	9.102	
•	.47532				.53958		1.719	.51241	6.466	**
Bromoform 4-Methyl-2-Pentanone	.12070		.10929	.10877	.10873	.11208	.784	.11260	4.323	
Toluene-d8 (SURR)		1,20804						1.22379		(Conc=50.0,50.0,5
Toluene	.91890		.76345	.74700	.76788	.76642	.830	.78669	8.291	*
2-Hexanone	.09059		.10052	.09388	.09390		.889	.09491	3.449	
Tetrachloroethene	.83533		.66857	.67711	.69011	.68342	.911	.70774	8.915	
Chlorobenzene		1.03579				-		1.05306	8.459	**
1,1,1,2-Tetrachloroethane	.63597	.54869	.56958	.58137	.59197	.58915	1.014	.58612	4.960	
Ethylbenzene		1.63410						1.70307	5.738	*
p & m-Xylene		1.42785						1.42529		(Conc=10.0,40.0,1
o-Xylene		1.38263						1.41441	8.160	(0010-1010)4010,1
Styrene	.75176	.63094	.63876	.63775	.62915	.62465	1.087	-65217	7.526	
Isopropylbenzene		1.76124						1.84215	6,199	
1,1,2,2-Tetrachloroethane	.71036		.61389	.60437	.58766	.57951	1.149	.61602	7.759	**
p-Bromofluorobenzene (SURR)	.64361	.64271	.65951	.65899	.65362	.64999	1.156	.65141		(Conc=50.0,50.0,5
1,2,3-Trichloropropane	.20940		.17330	.17068	. 16455	.16481	1.165	.17550	9.665	,,
Bromobenzene	.72229	.60313	.61126	.61394	.59938	.58648	1.172	.62275	7.985	
n-Propylbenzene	.50182	.43146	.42907	.43350	.42210	.41619	1.189	.43902	7.158	
2-Chlorotoluene	. 49854	.42504	.38566	.41990	.40835	.40089	1,199	.42306	9.346	
4-Chlorotoluene	. 48494	.41451	.45157	.40504	.40662	.39355	1.207	.42604	8.217	
1,3,5-Trimethylbenzene	1.82511	1.53462	1.55580	1.57308	1.55146	1.51298	1.211	1.59218	7.281	
tert-Butylbenzene	1.81233	1.50737	1.56450	1.53943	1.50263	1.46713	1.257	1.56556	8.009	
1,2,4-Trimethylbenzene	1.74271	1.47971	1.52558	1.51949	1.48664	1.45452	1.259	1.53478	6.855	
sec-Butylbenzene	2.40771	2.09909						2.15602	5.927	
1,3-Dichlorobenzene	1.21146		1.00100		.97563		1.298	1.01810	9.542	
p-Isopropyltoluene		1.67026						1.76029	6.908	
1,4-Dichlorobenzene	1.29679	1.07605				.97830		1.08770	10.171	
1,2-Dichtorobenzene	1.13388		.95637	.95775	.90765	.87035	1.356	. 9 6147	9.454	
n-Butylbenzene		1.77486						1.82740	6.263	
1,2-Dibromo-3-Chloropropane	. 13289		.14428		. 13656		1.457	. 13668	4.332	
1,2,4-Trichlorobenzene	.93731	.84305	.88679	.87259	.81863	.81125	1.582	.86160	5.496	
Naphthal ene		1.04615						1.11149	5.209	
Hexachlorobutadiene -	.90005	. 78090	.80922	.80459	.74905	.74878	1.614	.79877	7.013	
1,2,3-Trichlorobenzene	. 88330	.77862	.77532	.77301	.72377	.71405	1.638	.77468	7.764	

RF - Response factor (Subscript is amount in ppb)

RRT - Average Relative Retention Time (RT Std/RT Istd)

RF - Average Response Factor
%RSD - Percent Relative Standard Deviation
CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.7 (Cont) Results of the Initial Calibration WA # 0-098 Pier Drum

Instrument ID: GCMSD-2(3034A12982) Calibration Date: 01/24/95

Minimum RF for SPCC is 0.30

Maximum % RSD for CCC is 30.0%

Laboratory 1D	: >V2536 RF	>V2538 RF	>V2539 RF	>V2540 RF	>V2541 RF	>V2542 RF				
Compound	5.00	20.00	50.00	100.00	150.00	200.00	RRT	RF	% RSD	CCC SPCC
Dichlorodifluoromethane	3.87059	3.33391	3.22838	3.19722	3.04054	3.03838	.238	3.28484	9.401	
Chloromethane	1.33100	1.23270	1.23554	1.24097	1.19235	1.18036	.264	1.23549	4.292	**
Vinyl Chloride	1.42724	1.20120	1.16458	1.15769	1.09937	1.06762	.279	1.18628	10.743	*
3romomethane	1.57720	1.27261	1.14402	.94713	.90314	.91798	.335	1.12701	23.487	
Chloroethane	. 95468	. 74135	.73814	.68707	.46469	.45004	.353	.67266	28.329	•
Trichlorofluoromethane	4.81920	4.06191	3.99260	3.93498	3.79961	3.73424	.406	4.05709	9.675	
Acetone	.53048	.59603	.57807	.57474	.53902	.45065	.501	.54483	9.619	
.1-Dichloroethene	1.65195	1.38084	1.41141	1.35827	1.26812	1.26250	.520	1.38885	10.247	*
Carbon Disulfide	4.11539	3.52345	3.60498	3.59803	3.40153	3.42256	.592	3.61099	7.239	
Methylene Chloride	1.63957	1.44970	1.40192	1.38185	1.30150	1.28287	.624	1.40957	9.144	
rans-1,2-Dichloroethene	1.74445	1.49310	1.51774	1.45893	1.40142	1.34972	.704	1.49423	9.172	
lethyl-tertiary-butylether	2.53025	2.88152	2.87261	2.82026	2.75450	2.79136	.708	2.77508	4.657	
,1-Dichloroethane	3.17823	2.83320	2.82486	2.74891	2.64835	2.61711	.801	2.80844	7.182	**
-Butanone	.06774	.11941	. 12425	. 13992	. 13311	. 12528	.918	.11828	21.812	
,2-Dichloropropane	2.11446	1.91858	1.99826	2.03052	1.99384	2.01711	.943	2.01213	3.152	
is-1.2-Dichloroethene	1.75590	1.56003	1.53963	1.48402	1.41349	1.38872	.942	1.52363	8.679	
hloroform	3.82738	3.32782	3.39219	3.27721	3.13932	3.11947	.983	3.34723	7.709	*
,1-Dichloropropene	2.53002	2.23838	2.25801	2.18515	2.08972	2.05856	1.115	2.22664	7.565	
,2-Dichloroethane		-	_	2.11327				2.15505	5.744	
,2-Dichloroethane-d4 (SURR)	1.80818	1.86688	1,90156	1.90782	1.88785	1.95463	1.133	1.88782	2.577	(Conc=50.0,5
,1,1-Trichloroethane	.67387	.71797	.70861	.72047	.73814	.74145	.861	.71675	3.411	
arbon Tetrachloride	.20594		.50860	.55424	.59366	.59796	.904	.49380	29.712	
lenzene	.88000		.86666	.84167	.85064	.83854	.927	.86139	2.471	
richloroethene	.45948		.43135	.41560	.42563	.42090	1.047	.42868	3.745	
,2-Dichloropropane	.40905		.40002	.39911	.39908	.34417	1.075	.39471	6.524	*
ibromomethane	.39922		.39833	.37448	.38006	.36744	1.107	.38934	4.740	
romodichloromethane	.58803		.70545	.70742	.71957	.72447	1.111	.69009	7.399	
is-1,3-Dichloropropene	.43794		.47384	.47337	.47791	.46995	1.228	.47041	3.674	
rans-1,3-Dichloropropene	.33367					.41409	1.326	.40744	8.996	
1.1.2-Trichloroethane	.27552		.31088	. 29303	.30326		1.349	.29985	4.680	

RF - Response Factor (Subscript is amount in ppb)

RRT - Average Relative Retention Time (RT Std/RT 1std)

RF - Average Response Factor

%RSD - Percent Relative Standard Deviation
CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.7 (Cont) Results of the Initial Calibration WA # 0-098 Pier Drum

Instrument ID: GCMSD-2(3034A12982)

Calibration Date: 01/24/95

Minimum RF for SPCC is 0.30

Maximum % RSD for CCC is 30.0%

1,3-Dichloropropane .48698 .55994 .53783 .51608 .50381 .48895 1.397 .51560 5.573 Dibromochloromethane .51198 .66684 .65562 .64283 .68161 .66387 1.428 .63713 9.831 1,2-Dibromoethane .42110 .52417 .52814 .51233 .49747 .46737 1.464 .49176 8.344 Bromoform .33180 .47152 .50099 .52041 .51945 .53195 1.720 .47935 15.711 4-Methyl-2-Pentanone .07407 .11316 .11208 .11135 .11662 .12167 .782 .10816 15.834	ecc spec
Dibromochloromethane .51198 .66684 .65562 .64283 .68161 .66387 1.428 .63713 9.831 1,2-Dibromoethane .42110 .52417 .52814 .51233 .49747 .46737 1.464 .49176 8.344 Bromoform .33180 .47152 .50099 .52041 .51945 .53195 1.720 .47935 15.711 4-Methyl-2-Pentanone .07407 .11316 .11208 .11135 .11662 .12167 .782 .10816 15.834 Toluene-d8 (SURR) 1.23921 1.22496 1.24097 1.20978 1.24578 1.27976 .820 1.24008 1.893	**
Dibromochloromethane .51198 .66684 .65562 .64283 .68161 .66387 1.428 .63713 9.831 1,2-Dibromoethane .42110 .52417 .52814 .51233 .49747 .46737 1.464 .49176 8.344 Bromoform .33180 .47152 .50099 .52041 .51945 .53195 1.720 .47935 15.711 4-Methyl-2-Pentanone .07407 .11316 .11208 .11135 .11662 .12167 .782 .10816 15.834 Toluene-d8 (SURR) 1.23921 1.22496 1.24097 1.20978 1.24578 1.27976 .820 1.24008 1.893	**
1,2-Dibromoethane .42110 .52417 .52814 .51233 .49747 .46737 1.464 .49176 8.344 Bromoform .33180 .47152 .50099 .52041 .51945 .53195 1.720 .47935 15.711 4-Methyl-2-Pentanone .07407 .11316 .11208 .11135 .11662 .12167 .782 .10816 15.834 Toluene-d8 (SURR) 1.23921 1.22496 1.24097 1.20978 1.24578 1.27976 .820 1.24008 1.893	**
Bromoform .33180 .47152 .50099 .52041 .51945 .53195 1.720 .47935 15.711 4-Methyl-2-Pentanone .07407 .11316 .11208 .11135 .11662 .12167 .782 .10816 15.834 Toluene-d8 (SURR) 1.23921 1.22496 1.24097 1.20978 1.24578 1.27976 .820 1.24008 1.893	**
4-Methyl-2-Pentanone .07407 .11316 .11208 .11135 .11662 .12167 .782 .10816 15.834 Toluene-d8 (SURR) 1.23921 1.22496 1.24097 1.20978 1.24578 1.27976 .820 1.24008 1.893	
Toluene-d8 (SURR) 1.23921 1.22496 1.24097 1.20978 1.24578 1.27976 .820 1.24008 1.893	•
	(Conc=50.0,50.0,5
	*
2-Hexanone .09703 .14617 .14668 .14184 .13896 .13372 .890 .13407 14,002	
Tetrachloroethene .77317 .72219 .72002 .67760 .69136 .69114 .910 .71258 4.842	
Chlorobenzene 1.10151 .99880 1.02258 .99645 1.01279 1.03562 1.005 1.02796 3.785	**
1,1,1,2-Tetrachloroethane .55292 .56851 .58045 .58317 .60006 .61089 1.014 .58267 3.592	
Ethylbenzene 1.81262 1.62657 1.74459 1.64482 1.70056 1.75736 1.022 1.71442 4.137	*
	(Conc=10.0,40.0,1
o-Xylene 1.46768 1.40468 1.41423 1.36579 1.36497 1.39559 1.084 1.40216 2.705	• • • • • • • • • • • • • • • • • • • •
Styrene .63468 .64445 .66019 .64504 .65262 .67085 1.087 .65131 1.972	
Isopropylbenzene 1.88142 1.82038 1.84225 1.81909 1.78755 1.84085 1.137 1.83192 1.709	
1,1,2,2-Tetrachloroethane .52736 .59468 .59124 .58592 .59069 .61871 1.149 .58477 5.201	**
	(Conc=50.0,50.0,5
1,2,3-Trichtoropropane .13842 .16273 .16964 .16608 .16922 .17714 1.165 .16387 8.150	
Bromobenzene .61122 .60022 .60645 .60337 .58758 .59426 1.172 .60052 1.422	
n-Propylbenzene .43923 .42987 .42881 .43114 .41808 .42385 1.189 .42850 1.664	
2-Chilorotoluene .42353 .41582 .41164 .40546 .38880 .39146 1.199 .40612 3.382	
4-Chlorotaluene .46994 .42800 .44217 .43031 .43036 .43808 1.208 .43981 3.574	
1,3,5-Trimethylbenzene 1.64832 1.59380 1.57789 1.58833 1.57327 1.59946 1.213 1.59684 1.693	
tert-Butylbenzene 1.63453 1.54839 1.52996 1.53339 1.48713 1.50598 1.257 1.53990 3.324	
1,2,4-Trimethylbenzene 1.62601 1.51666 1.53400 1.54452 1.51601 1.52443 1.260 1.54360 2.708	
sec-Butylbenzene 2.25109 2.11870 2.15763 2.12733 2.10226 2.12359 1.286 2.14677 2.524	
1,3-Dichlorobenzene 1.05230 .97434 .96809 .97941 .93812 .93635 1.298 .97477 4.327	
p-Isopropyltoluene 1.80251 1.71410 1.71287 1.77950 1.72413 1.74507 1.308 1.74636 2.129	
1,4-Dichlorobenzene 1.10291 1.05704 1.05688 1.04998 1.00924 1.00552 1.313 1.04693 3.439	
1,2-Dichlorobenzene .91241 .94730 .90935 .92280 .88469 .89748 1.356 .91234 2.364	
n-Butylbenzene 1.83297 1.79140 1.82679 1.84606 1.79547 1.81464 1.362 1.81789 1.184	
1,2-Dibromo-3-Chloropropane .08863 .12806 .13799 .14184 .14247 .15083 1.457 .13164 16.960	
1,2,4-Trichtorobenzene .81189 .83411 .85700 .88606 .85526 .85434 1.582 .84978 2.930	
Naphthalene .78658 1.01931 1.07544 1.12158 1.10142 1.17141 1.604 1.04596 13.066	
Hexachlorobutadiene .77730 .77380 .79260 .82279 .77766 .77930 1.614 .78724 2.361	
1,2,3-Trichlorobenzene .72841 .76771 .74646 .77570 .76195 .75413 1.639 .75573 2.227	

RF - Response Factor (Subscript is amount in ppb)

RRT - Average Relative Retention Time (RT Std/RT Istd)

RF - Average Response Factor

%RSD - Percent Relative Standard Deviation CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)

098\DEL\AR\9503\QAQCVOA

0 114

Table 2.8 Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/13/95

Time: 13:05

Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 11/22/94

Minimum RF for SPCC is 0.30 Maximum % Diff for CCC is 25.0%

Compound	RF	RF	%Diff	ccc	SPCC
Dichlorodifluoromethane	2.75656	2.55358	7.36		
Chloromethane	1.10099	.95887	12.91		**
/inyl Chloride	1.06518	1.01573	4.64	*	
Bromomethane	1.00911	1.14206	13.17		
Chloroethane	.60516	.69096	14.18		
richlorofluoromethane	3.68991	3.88220	5.21		
Acetone	. 28846	.42185	46.24		
,1-Dichloroethene	1.31810	1.34489	2.03	*	
Carbon Disulfide	3.09811	3.21882	3.90		
ethylene Chloride	1.38705	1.31731	5.03		
rans-1,2-Dichloroethene	1.42359	1,47443	3.57		
ethyl-tertiary-butylether	2.76397	2.72785	1.31		
,1-Dichloroethane	2.74346	2.64074	3.74		**
-Butanone	.08201	.09353	14.04		
,2-Dichloropropane	2.03221	2.12645	4.64		
is-1,2-Dichloroethene	1.50057	1.53385	2.22		
hloroform	3.27645	3.27838	.06	*	
,1-Dichloropropene	2.03304	2.05881	1.27		
,2-Dichloroethane	2.12352	1.99255	6.17		
,2-Dichloroethane-d4 (SURR)	1.84429	1.71877	6.81		
,1,1-Trichloroethane	.70440	.71497	1.50		
arbon Tetrachloride	.49450	.48397	2.13		
enzene	.86300	.81681	5.35		
richloroethene	.44779	.43952	1.85		
.2-Dichloropropane	.41843	.34781	16.88	*	
ibromomethane	.42245	.38732	8.32		
romodichloromethane	.69864	.65445	6.33		
is-1,3-Dichloropropene	.48311	.47488	1.70		(Conc=50.
rans-1,3-Dichloropropene	.41905	.40980	2.21		(Conc=50.
,1,2-Trichloroethane	.32000	.28058	12.32		
,3-Dichloropropane	.53853	.50223	6.74		
ibromochloromethane	.64683	.60343	6.71		

- Response factor from daily standard file at 50.00 ppb RF

RF - Average Response Factor from Initial Calibration Form VI

%Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/13/95

Time: 13:05

Instrument ID: GCMSD-2(3034A12982)

Initial Calibration Date: 11/22/94

Minimum RF for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

Compound	RF	RF	X Diff	ccc	SPCC	
1,2-Dibromoethane	.52909	.50275	4.98			
Bromoform	.51241		4.76		**	
4-Methyl-2-Pentanone	.11260		6,22			
Toluene-d8 (SURR)		1.19818	2.09			
Toluene	.78669		2.60	*		
2-Hexanone	.09491	.12083	27.32			
Tetrachloroethene	.70774	.73895	4.41			
Chlorobenzene	1.05306	1.04191	1.06		**	
1,1,1,2-Tetrachloroethane	.58612	.58731	.20			
Ethylbenzene	1.70307	1.66329	2.34	*		
p & m-Xylene	1.42529	1.46331	2.67			(Conc=100.00)
o-Xylene	1.41441	1.39627	1.28			
Styrene	.65217	.63901	2.02			
Isopropylbenzene	1.84215	1.87564	1.82			
1,1,2,2-Tetrachloroethane	.61602	.54043	12.27		**	
p-Bromofluorobenzene (SURR)	.65141	.67324	3.35			
1,2,3-Trichloropropane	.17550	. 15655	10.80			
Bromobenzene	.62275	.63117	1.35			
n-Propylbenzene	.43902	.45095	2.72			
2-Chlorotoluene	.42306	.40535	4.19			
4-Chlorotoluene	.42604	.44780	5.11			
1,3,5-Trimethylbenzene	1.59218	1.59993	.49			
tert-Butylbenzene	1.56556	1.61927	3.43			
1,2,4-Trimethylbenzene	1.53478	1.54495	.66			
sec-Butylbenzene	2.15602		1.85			
1,3-Dichlorobenzene	1.01810		.65			
p-Isopropyltoluene	1.76029		3.67			
1,4-Dichlorobenzene	1.08770		3.78			
1,2-Dichlorobenzene	.96147		.71			
n-Butylbenzene		1.81106	.89			
1,2-Dibromo-3-Chloropropane		. 13300	2.69			
1,2,4-Trichlorobenzene	-	. 92075	6.86			
Naphthalene	1.11149		8.33			
Hexachlorobutadiene		.89094	11.54			
1,2,3-Trichlorobenzene	.77468	.78860	1.80			

- Response Factor from daily standard file at 50.00 ppb

RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/14/95

Time: 11:42

Initial Calibration Date: 11/22/94 Instrument ID: GCMSD-2(3034A12982)

Minimum RF for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

Compound	RF	RF	Diff	ccc	SPCC
Dichlorodifluoromethane	2.75656	2.50463	9.14		
Chloromethane	1.10099	1.04281	5.28		**
Vinyl Chloride	1.06518	1.03315	3.01	*	
Bromomethane	1.00911	1.13479	12.45		
Chloroethane	.60516	.68473	13.15		
Trichlorofluoromethane	3.68991	3.82987			
Acetone	.28846	.37413	29.70		
1,1-Dichloroethene	1.31810	1.35567	2.85	*	
Carbon Disulfide	3.09811	3.24807	4.84		
Methylene Chloride	1.38705	1.40269	1.13		
trans-1,2-Dichloroethene	1.42359	1.48376	4.23		
Methyl-tertiary-butylether	2. 7639 7	2.46775	10.72		
1,1-Dichloroethane	2.74346	2.72246	.77		**
2-Butanone	.08201	.08049	1.85		
2,2-Dichloropropane	2.03221	2.18718			
cis-1,2-Dichloroethene	1.50057	1.53246	2.13		
Chloroform	3.27645	3.27639	.00	*	
1,1-Dichloropropene	2.03304	2.11598	4.08		
1,2-Dichloroethane	2.12352	1.97882			
1,2-Dichloroethane-d4 (SURR)	1.84429		6.76		
1,1,1-Trichloroethane	. 7044 0	. 70398	.06		
Carbon Tetrachloride	.49450	.56042			
Benzene	.86300				
Trichloroethene	.44779		2.45		
1,2-Dichloropropane	.41843			*	
Dibromomethane	.42245	.35463	16.06		
Bromodichloromethane	.69864	.63877	8.57		
cis-1,3-Dichloropropene	. 48311	.45684	5.44		(Conc
trans-1,3-Dichloropropene	.41905				(Conc
1,1,2-Trichloroethane	.32000				
1,3-Dichloropropane	.53853		9.31		
Dibromochloromethane	. 64683	.60466	6.52		

RF - Response Factor from daily standard file at 50.00 ppb

RF - Average Response Factor from Initial Calibration Form VI

%Dirf - % Difference from original average or curve

CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/14/95

Time: 11:42

Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 11/22/94

Minimum RF for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

Compound	RF	RF	%Diff	CCC	SPCC	
1,2-Dibromoethane	.52909	.48850	7.67			
Bromoform	.51241		8.11		**	
4-Methyl-2-Pentanone	.11260		20.95			
Toluene-d8 (SURR)		1.22117	.21			
Toluene	.78669	.77357	1.67	*		
2-Hexanone	.09491	.10288	8.40			
Tetrachloroethene	.70774	.73834	4.32			
Chlorobenzene	1.05306	1.04481	.78		**	
1,1,1,2-Tetrachloroethane	.58612	.58108	.86			
Ethylbenzene	1.70307	1.65794	2.65	*		
p & m-Xylene	1.42529	1.48248	4.01			(Conc=100.00)
o-Xylene	1.41441	1.40737	.50			
Styrene	.65217	.64200	1.56			
Isopropylbenzene	1.84215	1.87787	1.94			
1,1,2,2-Tetrachloroethane	.61602	.51245	16.81		**	
p-Bromofluorobenzene (SURR)	.65141	.67719	3.96			
1,2,3-Trichloropropane	. 17550	. 15249	13.11			
Bromobenzene	.62275	.62966	1.11			
n-Propylbenzene	.43902		4.07			
2-Chlorotoluene	. 42306	.40451	4.39			
4-Chlorotoluene	. 42604	.47032	10.39			
1,3,5-Trimethylbenzene	1.59218		.69			•
tert-Butylbenzene	1.56556		4.56			
1,2,4-Trimethylbenzene	1.53478		1.36			
sec-Butylbenzene	2.15602		2.00			
1,3-Dichlorobenzene	1.01810		.97			
p-Isopropyltoluene	1.76029		4.23			
1,4-Dichlorobenzene	1.08770		2.98			
1,2-Dichlorobenzene	.96147	.95389	.79			
n-Butylbenzene	1.82740		. 19			
1,2-Dibromo-3-Chloropropane	. 13668		11.04			
1,2,4-Trichlorobenzene	.86160		6.88			
Naphthalene	1,11149		12.06			
Hexachlorobutadiene	.79877		13.85			
1,2,3-Trichlorobenzene	.77468	.77641	.22			

- Response Factor from daily standard file at 50.00 ppb

RF - Average Response factor from Initial Calibration Form VI %Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/17/95

Time: 14:32 Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 11/22/94

Minimum $\overline{\text{RF}}$ for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

Compound	RF	RF	%Diff	ССС	SPCC
Dichlorodifluoromethane	2.75656	2.49224	9.59		
Chloromethane	1.10099	1.01824	7.52		**
/inyl Chloride	1.06518	1.07157	.60	*	
3romomethane	1.00911	1.24400	23.28		
Chloroethane	.60516	.73023	20.67		
Trichlorofluoromethane	3.68991	3.91212	6.02		
lcetone	. 28846	.44064	52.75		
1,1-Dichloroethene	1.31810	1.40152	6.33	*	
Carbon Disulfide	3.09811	3.34504	7.97		
Methylene Chloride	1.38705	1.42499	2.74		
rans-1,2-Dichloroethene	1.42359	1.53474	7.81		
Methyl-tertiary-butylether	2.76397	2.64108	4.45		
,1-Dichloroethane	2.74346	2.79274	1.80		**
?-Butanone	.08201	.09665	17.85		
2,2-Dichloropropane	2.03221	2.22924	9.70		
is-1,2-Dichloroethene	1.50057	1.58243	5.46		
Chloroform	3.27645	3.40136	3.81	*	
,1-Dichloropropene	2.03304	2.09723	3.16		
,2-Dichloroethane	2.12352	2.10906	.68		
,2-Dichloroethane-d4 (SURR)	1.84429	1.75566	4.81		
,1,1-Trichloroethane	.70440	.73263	4.01		
arbon Tetrachloride	.49450	.50930	2.99		
lenzene	.86300	.85655	.75		
richloroethene	.44779	.43745			
,2-Dichloropropane	.41843	.40573		*	
ibromomethane	.42245	.42741	1.17		
romodichloromethane	.69864	.70826	1.38		
is-1,3-Dichloropropene	.48311	.50675	4.89		(Conc=50.0
rans-1,3-Dichloropropene	.41905				(Conc≈50.0
,1,2-Trichloroethane	.32000	.30797	3.76		
,3-Dichloropropane	.53853	.53966	.21		
ibromochloromethane	.64683	.71826	11.04		

RF - Response Factor from daily standard file at 50.00 ppb

- Average Response Factor from Initial Calibration Form VI

%Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/17/95

Time: 14:32

Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 11/22/94

Minimum RF for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

1,2-Dibromoethane	Compound	RF	RF	X Diff	CCC	SPCC	
## ## ## ## ## ## ## ## ## ## ## ## ##	1.2-Dibromoethane	.52909	.53886	1.85			
Toluene		.51241	.53444	4.30		**	
Toluene	4-Methyl-2-Pentanone	.11260	. 10554	6.28			
Toluene			1.19695	2.19			
Tetrachloroethene Chlorobenzene 1.05306 1.07613 2.19 ** 1,1,2-Tetrachloroethane Ethylbenzene 1.70307 1.76439 3.60 * p & m-xylene 1.42529 1.48963 4.51 (Conc=100.00) o-Xylene 1.41441 1.45118 2.60 Styrene 1.5217 .66155 1.44 Isopropylbenzene 1.84215 1.94756 5.72 1,1,2,2-Tetrachloroethane p-Bromofluorobenzene (SURR) 1.2,3-Trichloropropane 1.7550 1.7123 2.44 Bromobenzene 62275 .65918 5.85 n-Propylbenzene 62275 .65918 5.85 n-Propylbenzene 42306 .40903 3.32 4-Chlorotoluene 42306 .40903 3.32 4-Chlorotoluene 42604 .49614 16.45 1,3,5-Trimethylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 1.56556 1.67700 7.12 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08170 1.14138 4.94 1,2-Dichlorobenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 1.4181 3.76 1,2,4-Trichlorobenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 1.4181 3.76 1,2,4-Trichlorobenzene 1.3668 1.4181 3.76 1.27 1.391484 14.53		.78669	.79232	.72	*		
Chlorobenzene 1.05306 1.07613 2.19 *** 1,1,1,2-Tetrachioroethane 58612 .60649 3.47 Ethylbenzene 1.70307 1.76439 3.60 * p & m-Xylene 1.42529 1.48963 4.51 (Conc=100.00) o-Xylene 1.41441 1.45118 2.60 Styrene .65217 .66155 1.44 Isopropylbenzene 1.84215 1.94756 5.72 1,1,2,2-Tetrachioroethane .61602 .58083 5.71 *** p-Bromofluorobenzene (SURR) .65141 .68763 5.56 1,2,3-Trichloropropane .17550 .17123 2.44 Bromobenzene .62275 .65918 5.85 n-Propylbenzene .42306 .40903 3.32 4-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42604 .49614 16.45 1,3,5-Trimethylbenzene 1.55218 1.64760 3.48 tetr-Butylbenzene 1.55556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	2-Hexanone	.09491	.11717	23.45			
1,1,1,2-Tetrachloroethane Ethylbenzene p & m-Xylene p & m-Xylene 1.42529 1.48963 4.51 o-Xylene 1.41441 1.45118 2.60 Styrene 1.55217 .66155 1.44 Isopropylbenzene 1.84215 1.94756 5.72 1,1,2,2-Tetrachloroethane p-Bromofluorobenzene (SURR) 1,2,3-Trichloropropane 1.7550 .17123 2.44 Bromobenzene 1.62275 .65918 5.85 n-Propylbenzene 1.43902 .46245 5.34 2-Chlorotoluene 1.42506 .40903 3.32 4-Chlorotoluene 1.59218 1.64760 3.48 tetr-Butylbenzene 1.59218 1.64760 3.48 tetr-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 1.153478 1.60743 4.73 sec-Butylbenzene 1.15002 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.2445 1.01712 5.79 n-Butylbenzene 1.86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene 1.79877 .91484 14.53	Tetrachloroethene	.70774	.75584	6.80			
Ethylbenzene 1.70307 1.76439 3.60 * p & m-Xylene 1.42529 1.48963 4.51 (Conc=100.00) o-Xylene 1.41441 1.45118 2.60 Styrene 6.65217 .66155 1.44 Isopropylbenzene 1.84215 1.94756 5.72 1,1,2,2-Tetrachloroethane 61602 .58083 5.71 ** p-Bromofluorobenzene (SURR) 65141 .68763 5.56 1,2,3-Trichloropropane 1.7550 .17123 2.44 Bromobenzene 62275 .65918 5.85 n-Propylbenzene 62275 .65918 5.85 n-Propylbenzene 643902 .46245 5.34 2-Chlorotoluene 642604 .49614 16.45 1,3,5-Trimethylbenzene 61.59218 1.64760 3.48 tert-Butylbenzene 61.53478 1.60743 4.73 sec-Butylbenzene 61.53478 1.60743 4.73 sec-Butylbenzene 61.01810 1.07923 6.00 p-Isopropyltoluene 61.76029 1.89192 7.48 1,4-Dichlorobenzene 61.08770 1.14138 4.94 1,2-Dichlorobenzene 61.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 61.0868 14181 3.76 1,2,4-Trichlorobenzene 61.11149 1.08289 2.57 Hexachlorobutadiene 79877 .91484 14.53	Chlorobenzene	1.05306	1.07613	2.19		**	
p & m-Xylene 1.42529 1.48963 4.51 (Conc=100.00) o-Xylene 1.41441 1.45118 2.60 Styrene .65217 .66155 1.44 Isopropylbenzene 1.84215 1.94756 5.72 1,1,2,2-Tetrachloroethane .61602 .58083 5.71 *** p-Bromofluorobenzene (SURR) .65141 .68763 5.56 1,2,3-Trichloropropane .17550 .17123 2.44 Bromobenzene .62275 .65918 5.85 n-Propylbenzene .62275 .65918 5.85 n-Propylbenzene .42306 .40903 3.32 4-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42604 .49614 16.45 1,3,5-Trimethylbenzene 1.55556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.8668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1	1,1,1,2-Tetrachloroethane	.58612	.60649	3.47			
o-Xylene 1.41441 1.45118 2.60 Styrene .65217 .66155 1.44 Isopropylbenzene 1.84215 1.94756 5.72 1,1,2,2-Tetrachloroethane .61602 .58083 5.71 *** p-Bromofluorobenzene (SURR) .65141 .68763 5.56 1,2,3-Trichloropropane .17550 .17123 2.44 Bromobenzene .62275 .65918 5.85 n-Propylbenzene .43902 .46245 5.34 2-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42604 .49614 16.45 1,3,5-Trimethylbenzene 1.59218 1.64760 3.48 tert-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dibromo-3-Chloropropane 1.82740 1.86937 2.30 1,2,4-Trichlorobenzene 1.8668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene 7.9877 .91484 14.53 <td>Ethylbenzene</td> <td>1.70307</td> <td>1.76439</td> <td>3.60</td> <td>*</td> <td></td> <td></td>	Ethylbenzene	1.70307	1.76439	3.60	*		
Styrene .65217 .66155 1.44 Isopropylbenzene 1.84215 1.94756 5.72 1,1,2,2-Tetrachloroethane .61602 .58083 5.71 *** p-Bromofluorobenzene (SURR) .65141 .68763 5.56 1,2,3-Trichloropropane .17550 .17123 2.44 Bromobenzene .62275 .65918 5.85 n-Propylbenzene .42306 .40903 3.32 4-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42504 .49614 16.45 1,3,5-Trimethylbenzene 1.59218 1.64760 3.48 tert-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 1.4181 3.76	p & m-Xylene	1.42529	1.48963	4.51			(Conc=100.00)
Isopropylbenzene 1.84215 1.94756 5.72 1,1,2,2-Tetrachloroethane .61602 .58083 5.71 *** p-Bromofluorobenzene (SURR) .65141 .68763 5.56 1,2,3-Trichloropropane .17550 .17123 2.44 Bromobenzene .62275 .65918 5.85 n-Propylbenzene .43902 .46245 5.34 2-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42504 .49614 16.45 1,3,5-Trimethylbenzene 1.59218 1.64760 3.48 tert-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 1.4181 3.76 1,2,4-Trichlorobenzene 86160 .96730 12.27 </td <td>o-Xylene</td> <td>1.41441</td> <td>1.45118</td> <td>2.60</td> <td></td> <td></td> <td></td>	o-Xylene	1.41441	1.45118	2.60			
1,1,2,2-Tetrachloroethane p-Bromofluorobenzene (SURR) 1,2,3-Trichloropropane 1,7550 1,12,3-Trichloropropane 1,7550 1,7123 2,44 Bromobenzene 1,62275 1,65918 1,85 1,85 1,87 1,87 1,87 1,87 1,87 1,87 1,87 1,87	Styrene	.65217	.66155	1.44			
p-Bromofluorobenzene (SURR)	Isopropylbenzene	1.84215	1.94756	5.72			
1,2,3-Trichloropropane .17550 .17123 2.44 Bromobenzene .62275 .65918 5.85 n-Propylbenzene .43902 .46245 5.34 2-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42604 .49614 16.45 1,3,5-Trimethylbenzene 1.59218 1.64760 3.48 tert-Butylbenzene 1.5556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 2.15602 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	1,1,2,2-Tetrachloroethane	.61602	.58083	5.71		**	
Bromobenzene .62275 .65918 5.85 n-Propylbenzene .43902 .46245 5.34 2-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42604 .49614 16.45 1,3,5-Trimethylbenzene 1.59218 1.64760 3.48 tert-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 2.15602 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	p-Bromofluorobenzene (SURR)	.65141	.68763	5.56			
n-Propylbenzene .43902 .46245 5.34 2-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42604 .49614 16.45 1,3,5-Trimethylbenzene 1.59218 1.64760 3.48 tert-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 2.15602 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	1,2,3-Trichloropropane	.17550	.17123	2.44			
2-Chlorotoluene .42306 .40903 3.32 4-Chlorotoluene .42604 .49614 16.45 1,3,5-Trimethylbenzene 1.59218 1.64760 3.48 tert-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 2.15602 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	Bromobenzene	.62275	.65918	5.85			
4-Chlorotoluene .42604 .49614 16.45 1,3,5-Trimethylbenzene 1.59218 1.64760 3.48 tert-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 2.15602 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	n-Propylbenzene	.43902	.46245	5.34			
1,3,5-Trimethylbenzene 1.59218 1.64760 3.48 tert-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 2.15602 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	2-Chlorotoluene	.42306	.40903	3.32			
tert-Butylbenzene 1.56556 1.67700 7.12 1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 2.15602 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	4-Chlorotoluene	.42604	.49614	16.45			
1,2,4-Trimethylbenzene 1.53478 1.60743 4.73 sec-Butylbenzene 2.15602 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	1,3,5-Trimethylbenzene	1.59218	1.64760				
sec-Butylbenzene 2.15602 2.26612 5.11 1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane 1.3668 14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	tert-Butylbenzene	1.56556	1.67700	7.12			
1,3-Dichlorobenzene 1.01810 1.07923 6.00 p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane .13668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	1,2,4-Trimethylbenzene	1.53478	1.60743				
p-Isopropyltoluene 1.76029 1.89192 7.48 1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane .13668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	sec-Butylbenzene						
1,4-Dichlorobenzene 1.08770 1.14138 4.94 1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane .13668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	1,3-Dichlorobenzene						
1,2-Dichlorobenzene .96147 1.01712 5.79 n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane .13668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	p-Isopropyltoluene			–			
n-Butylbenzene 1.82740 1.86937 2.30 1,2-Dibromo-3-Chloropropane .13668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	1,4-Dichlorobenzene	1.08770	1.14138				
1,2-Dibromo-3-Chloropropane .13668 .14181 3.76 1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	1,2-Dichlorobenzene						
1,2,4-Trichlorobenzene .86160 .96730 12.27 Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53	· · · · · · · · · · · · · · · · · · ·						
Naphthalene 1.11149 1.08289 2.57 Hexachlorobutadiene .79877 .91484 14.53							
Hexachlorobutadiene .79877 .91484 14.53	1,2,4-Trichlorobenzene						
1,2,3-Trichlorobenzene .77468 .84018 8.46			-				
	1,2,3-Trichlorobenzene	.77468	.84018	8.46			

RF - Response Factor from daily standard file at 50.00 ppb

RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/18/95

Time: 08:27

Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 01/18/95

Minimum RF for SPCC is 0.30 Maximum % Diff for CCC is 25.0%

	Compound	RF	RF	X Diff	ccc	SPCC
Vinyl Chloride 1.03762 1.03141 .60 * Bromomethane 1.07394 1.00215 6.68 * Chloroethane .61678 .70076 13.62 * Trichlorofluoromethane 3.77576 3.72338 1.39 Acetone .46918 .47822 1.93 1,1-Dichloroethene 1.34780 1.33850 .69 * Carbon Disulfide 3.16441 3.07738 2.75 * Methylene Chloride 1.39021 1.37674 .97 * trans-1,2-Dichloroethene 1.48094 1.47818 .19 * Methyl-tertiary-butylether 2.75552 2.77875 .84 * * 1,1-Dichloroethane 2.075552 2.77875 .84 * * 2,-Butanone 2.10782 2.10790 .00 * * * 2,-Dichloropropene 2.10782 2.10790 .00 * * * 1,1-Dichloropropene 2.07881 2.05973	Dichlorodifluoromethane	2,41009	2.38801	.92		
Bromomethane	Chloromethane	1.04526	1.00295	4.05		**
Chloroethane .61678 .70076 13.62 Trichlorofluoromethane 3.77576 3.72338 1.39 Acetone .46918 .47822 1.93 1,1-Dichloroethene 1.34780 1.33850 .69 * Carbon Disulfide 3.16441 3.07738 2.75 Methylene Chloride 1.39021 1.37674 .97 trans-1,2-Dichloroethene 1.48094 1.47818 .19 Methyl-tertiary-butylether 2.75552 2.77875 .84 1,1-Dichloroethane 2.72588 2.72658 .03 ** 2-Butanone 1.0054 .10345 2.89 2,2-Dichloroethane 1.053397 1.55992 1.69 Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92 1,2-Dichloroethane 2.05552 2.04742 .39 1,2-Dichloroethane 2.05552 2.04742 .39 1,1,1-Trichloroethane .69995 .71418 2.03 Carbon Tetrachloride .84339 .85451 1.32 </td <td>Vinyl Chloride</td> <td>1.03762</td> <td>1.03141</td> <td>.60</td> <td>*</td> <td></td>	Vinyl Chloride	1.03762	1.03141	.60	*	
Trichlorofluoromethane	Bromomethane	1.07394	1.00215	6.68		
Acetone	Chloroethane	.61678	.70076	13.62		
1,1-Dichloroethene 1.34780 1.33350 .69 Carbon Disulfide 3.16441 3.07738 2.75 Methylene Chloride 1.39021 1.37674 .97 trans-1,2-Dichloroethene 1.48094 1.47818 .19 Methyl-tertiary-butylether 2.75552 2.77875 .84 1,1-Dichloroethane 2.72588 2.72658 .03 *** 2-Butanone .10054 .10345 2.89 2,2-Dichloropropane 2.10782 2.10790 .00 cis-1,2-Dichloroethene 1.53397 1.55992 1.69 Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92 1.2-Dichloroethane 2.05552 2.04742 .39 1.2-Dichloroethane .69995 .71418 2.03 <td< td=""><td>Trichlorofluoromethane</td><td>3.77576</td><td>3.72338</td><td>1.39</td><td></td><td></td></td<>	Trichlorofluoromethane	3.77576	3.72338	1.39		
Carbon Disulfide 3.16441 3.07738 2.75 Methylene Chloride 1.39021 1.37674 .97 trans-1,2-Dichloroethene 1.48094 1.47818 .19 Methyl-tertiary-butylether 2.75552 2.77875 .84 1,1-Dichloroethane 2.72588 2.72658 .03 ** 2-Butanone .10054 .10345 2.89 2,2-Dichloropropane 2.10782 2.10790 .00 cis-1,2-Dichloroethene 1.53397 1.55992 1.69 Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92 1,2-Dichloroethane 2.05552 2.04742 .39 1,2-Dichloroethane .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .66356 2.01 66356	Acetone	.46918	.47822	1.93		
Methylene Chloride 1.39021 1.37674 .97 trans-1,2-Dichloroethene 1.48094 1.47818 .19 Methyl-tertiary-butylether 2.75552 2.77875 .84 1,1-Dichloroethane 2.72588 2.72658 .03 ** 2-Butanone .10054 .10345 2.89 2,2-Dichloropropane 2.10782 2.10790 .00 cis-1,2-Dichloroethene 1.53397 1.55992 1.69 Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92 1.2-Dichloroethane 2.07881 2.05973 .92 1,2-Dichloroethane 2.05552 2.04742 .39 1.70 1.1,1-Trichloroethane .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 8 Benzene .84339 .85451 1.32 Trichloroethane .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .66356 2.01 2.01	1,1-Dichloroethene	1.34780	1.33850	.69	*	
trans-1,2-Dichloroethene 1.48094 1.47818 .19 Methyl-tertiary-butylether 2.75552 2.77875 .84 1,1-Dichloroethane 2.72588 2.72658 .03 ** 2-Butanone .10054 .10345 2.89 2,2-Dichloropropane 2.10782 2.10790 .00 cis-1,2-Dichloroethene 1.53397 1.55992 1.69 Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92 1,2-Dichloroethane 2.05552 2.04742 .39 1,2-Dichloroethane 2.05552 2.04742 .39 1,2-Dichloroethane .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .667719 .66356 2.01 * cis-1,3-Dichloropropene .47831 .49244	Carbon Disulfide	3.16441	3.07738	2.75		
Methyl-tertiary-butylether 2.75552 2.77675 .84 1,1-Dichloroethane 2.72588 2.72658 .03 ** 2-Butanone .10054 .10345 2.89 2,2-Dichloropropane 2.10782 2.10790 .00 cis-1,2-Dichloroethane 1.53397 1.55992 1.69 Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92	Methylene Chloride	1.39021	1.37674	.97		
1,1-Dichloroethane 2.72588 2.72658 .03 *** 2-Butanone .10054 .10345 2.89 2,2-Dichloropropane 2.10782 2.10790 .00 cis-1,2-Dichloroethene 1.53397 1.55992 1.69 Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92 1,2-Dichloroethane 2.05552 2.04742 .39 .92 1,2-Dichloroethane .69995 .71418 2.03 Carbon Tetrachloride .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .66356 2.01 .01 .01 .01 .02 .01 .02 .03 .03 .03 .03 .03 .03 .03 .03 .03 .03 .03 .03 .03 <td< td=""><td>trans-1,2-Dichloroethene</td><td>1.48094</td><td>1.47818</td><td>. 19</td><td></td><td></td></td<>	trans-1,2-Dichloroethene	1.48094	1.47818	. 19		
2-Butanone	Methyl-tertiary-butylether	2.75552	2.77875	.84		
2,2-Dichloropropane 2.10782 2.10790 .00 cis-1,2-Dichloroethene 1.53397 1.55992 1.69 Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92 1,2-Dichloroethane 2.05552 2.04742 .39 1,2-Dichloroethane-d4 (SURR) 1.78628 1.75591 1.70 1,1,1-Trichloroethane .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .41354 .38513 6.87 Bromodichloromethane .667719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	1,1-Dichloroethane	2.72588	2.72658	. 03		**
cis-1,2-Dichloroethene 1.53397 1.55992 1.69 Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92 1,2-Dichloroethane 2.05552 2.04742 .39 1,2-Dichloroethane .6995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .41354 .38513 6.87 Bromodichloromethane .67719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	2-Butanone	. 10054	. 10345	2.89		
Chloroform 3.27541 3.30868 1.02 * 1,1-Dichloropropene 2.07881 2.05973 .92 1,2-Dichloroethane 2.05552 2.04742 .39 1,2-Dichloroethane-d4 (SURR) 1.78628 1.75591 1.70 1,1,1-Trichloroethane .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .41354 .38513 6.87 Bromodichloromethane .67719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	2,2-Dichloropropane	2.10782	2.10790	.00		
1,1-Dichloropropene 2.07881 2.05973 .92 1,2-Dichloroethane 2.05552 2.04742 .39 1,2-Dichloroethane-d4 (SURR) 1.78628 1.75591 1.70 1,1,1-Trichloroethane .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 Pibromomethane .41354 .38513 6.87 Bromodichloromethane .667719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	cis-1,2-Dichloroethene	1.53397	1.55992			
1,2-Dichloroethane 2.05552 2.04742 .39 1,2-Dichloroethane-d4 (SURR) 1.78628 1.75591 1.70 1,1,1-Trichloroethane .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 Pibromomethane .41354 .38513 6.87 Bromodichloromethane .667719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	Chloroform	3.27541	3.30868	1.02	*	
1,2-Dichloroethane-d4 (SURR) 1.78628 1.75591 1.70 1,1,1-Trichloroethane .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .41354 .38513 6.87 Bromodichloromethane .667719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	1,1-Dichloropropene	2.07881	2.05973	.92		
1,1,1-Trichloroethane .69995 .71418 2.03 Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .41354 .38513 6.87 Bromodichloromethane .67719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	1,2-Dichloroethane					
Carbon Tetrachloride .50468 .47367 6.14 Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .41354 .38513 6.87 Bromodichloromethane .67719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	1,2-Dichloroethane-d4 (SURR)	-				
Benzene .84339 .85451 1.32 Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .41354 .38513 6.87 Bromodichloromethane .67719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	1,1,1-Trichloroethane					
Trichloroethene .44451 .45354 2.03 1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .41354 .38513 6.87 Bromodichloromethane .67719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	Carbon Tetrachloride					
1,2-Dichloropropane .37881 .33551 11.43 * Dibromomethane .41354 .38513 6.87 Bromodichloromethane .67719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	Benzene	. 84339	. 85451			
Dibromomethane .41354 .38513 6.87 Bromodichloromethane .67719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.00) trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.00) 1,1,2-Trichloroethane .30073 .30322 .83	Trichloroethene	.44451	.45354	2.03		
Bromodichloromethane .67719 .66356 2.01 cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	1,2-Dichloropropane	.37881	. 33551		*	
cis-1,3-Dichloropropene .47831 .49244 2.95 (Conc=50.0 trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 trans-1,1,2-Trichloropropene .30073 .30322 .83	Dibromomethane					
trans-1,3-Dichloropropene .40463 .42874 5.96 (Conc=50.0 1,1,2-Trichloroethane .30073 .30322 .83	Bromodichloromethane	.67719				
1,1,2-Trichloroethane .30073 .30322 .83	cis-1,3-Dichloropropene					(Conc=50.00)
, , -	trans-1,3-Dichloropropene					(Conc=50.00)
1,3-Dichloropropane .50907 .52452 3.04	1,1,2-Trichloroethane					
Dibromochloromethane .65430 .67871 3.73	Dibromochloromethane	.65430	.67871	3.73		

RF - Response Factor from daily standard file at 50.00 ppb

RF - Average Response Factor from Initial Calibration Form VI

%Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/18/95

Time: 08:27

Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 01/18/95

Minimum RF for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

1,2-Dibromoethane	Compound	RF	RF	% Diff	CCC	SPCC	
## ## ## ## ## ## ## ## ## ## ## ## ##	1 2-Dibromosthans	50051	52000	7 R7			
## ## ## ## ## ## ## ## ## ## ## ## ##						**	
Toluene-d8 (SURR) 1.20579 1.20925 .29 Toluene .78752 .79244 .63 * 2-Hexanone .12437 .13527 8.76 Tetrachloroethene .74202 .75230 1.38 Chlorobenzene 1.05861 1.08336 2.34 ** 1,1,2-Tetrachloroethane .58493 .59720 2.10 Ethylbenzene 1.74668 1.74948 1.6 * p & m-Xylene 1.41406 1.47559 4.35 (Conc=100.00) o-Xylene 1.43177 1.44755 1.10 Styrene .64562 .65686 1.74 Isopropylbenzene 1.89480 1.93251 1.99 1,1,2,2-Tetrachloroethane p-Bromofluorobenzene (SURR) 1.67535 .69319 2.64 1,2,3-Trichloropropane .63774 .65675 2.98 n-Propylbenzene .63774 .65675 2.98 n-Propylbenzene .45524 .46106 1.28 2-Chlorotoluene .45803 .43718 4.55 1,3,5-Trimethylbenzene 1.62309 1.61498 .50 tetr-Butylbenzene 1.63687 1.57045 .09 sec-Butylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.8422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.3427 1.3950 3.89 1,2,4-Trichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.8422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.13427 1.3950 3.89 1,2,3-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.10858 .85528 .73 1,2,3-Trichlorobenzene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	4						
Toluene	•						
2-Mexanone Tetrachloroethene Tetrachloroethene Thirdenic 1.05861 1.08336 2.34	The state of the s				*		
Tetrachloroethene Chlorobenzene 1.05861 1.08336 2.34 ** 1.1,1,2-Tetrachloroethane Ethylbenzene p & m-Xylene 0.4568 1.74948 .16 * p & m-Xylene 1.41406 1.47559 4.35 (Conc=100.00) 0-Xylene 1.43177 1.44755 1.10 Styrene 1.43177 1.44755 1.10 Styrene 1.4568 1.93251 1.99 1.1,2,2-Tetrachloroethane p-Bromofluorobenzene (SURR) 1.2,3-Trichloropropane 1.6768 .17259 2.93 Bromobenzene 1.63774 .65675 2.98 n-Propylbenzene 1.63774 .65675 2.98 n-Propylbenzene 2.616764 .44855 7.40 4-Chlorotoluene 4.5803 .43718 4.55 1.3,5-Trimethylbenzene 1.62309 1.61498 .50 tert-Butylbenzene 1.63697 1.57045 .09 sec-Butylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 1.03711 1.06074 2.28 p-lsopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.03711 1.06074 2.28 p-lsopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.21822 1.12822 .18 1,2-Dichlorobenzene 1.2622 1.12822 .18 1,2-Dichlorobenzene 1.2624 1.282916 .82 1,2-Dibromo-3-Chloropropane 1.2,4-Trichlorobenzene 1.2938 1.10858 1.84 Hexachlorobutadiene 1.2938 1.10858 1.84 Hexachlorobutadiene 1.80946 .79040 2.35	,						
Chlorobenzene 1.05861 1.08336 2.34 *** 1,1,1,2-Tetrachloroethane .58493 .59720 2.10 Ethylbenzene 1.74668 1.74948 .16 * p & m-Xylene 1.44106 1.47559 4.35 (Conc=100.00) o-Xylene 1.43177 1.44755 1.10 (Conc=100.00) Styrene .64562 .65686 1.74 1.50 (Conc=100.00) 1,2,2-Tetrachloroethane .57686 .59682 3.46 ** p-Bromofluorobenzene (SURR) .67535 .69319 2.64 ** 1,2,3-Trichloropropane .16768 .17259 2.93 ** Bromobenzene .63774 .65675 2.98 ** n-Propylbenzene .45524 .46106 1.28 2-Chlorotoluene .41764 .44855 7.40 4-Chlorotoluene .45803 .43718 4.55 1,3,5-Trimethylbenzene 1.61365 1.63576 1.37 1,2,4-Trimethylbenzene 1.6365 1.63576 1.37 1,2,4-Trimethylbenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene .98178 .98022 .16<							
1,1,1,2-Tetrachloroethane Ethylbenzene Ethylbenzene p & m-Xylene 0-Xylene 1.441406 1.47559 0-Xylene 1.43177 1.44755 1.10 Styrene 1.43177 1.44755 1.10 Styrene 1.43177 1.44755 1.10 Styrene 1.43177 1.44755 1.10 Styrene 1.89480 1.93251 1.99 1,1,2,2-Tetrachloroethane p-Bromofluorobenzene (SURR) 1,2,3-Trichloropropane 1.6783 1.7259 2.93 Bromobenzene 1.63774 6.5675 2.98 n-Propylbenzene 1.63774 6.5675 2.98 n-Propylbenzene 1.45524 4.6106 1.28 2-Chlorotoluene 4.41764 4.4855 7.40 4-Chlorotoluene 4.41764 4.4855 7.40 4-Chlorotoluene 1.63309 1.61498 .50 tert-Butylbenzene 1.63309 1.61498 .50 tert-Butylbenzene 1.63576 1.37 1,2,4-Trimethylbenzene 1.58897 1.577045 .09 sec-Butylbenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.12938 1.10858 1.84 Hexachlorobutadiene 1.12938 1.10858 1.84 Hexachlorobenzene 1.80946 .79040 2.35	:					**	
Ethylbenzene p & m-Xylene p & m-Xylene o-Xylene 1.41406 1.47559 1.10 Styrene 1.43177 1.44755 1.10 Styrene 1.89480 1.93251 1.99 1,1,2,2-Tetrachloroethane p-Bromofluorobenzene (SURR) 1,2,3-Trichloropropane 1.6768 1.7259 2.93 Bromobenzene 1.63774 .65675 2.98 n-Propylbenzene 2-Chlorotoluene 4.4552 4.46106 4-Chlorotoluene 4.4764 4.4855 7.40 4-Chlorotoluene 4.4764 4.4855 7.40 4-Chlorotoluene 1.62309 1.61498 .50 tert-Butylbenzene 1.61365 1.63576 1.37 1,2,4-Trimethylbenzene 1.61365 1.63576 1.37 1,2,4-Trimethylbenzene 1.80992 1.83815 1,3-Dichlorobenzene 1.80992 1.83815 1,4-Dichlorobenzene 1.12622 1.12822 1.8 1,2-Dichlorobenzene 1.84422 1.82916 1.2-Dibromo-3-Chloropropane 1.12938 1.10858 1.84 Hexachlorobutadiene 1.12938 1.10858 1.84 Hexachlorobutadiene 1.80966 .79040 2.35 (Conc=100.00) ** (Conc=100.00) (Conc=100.00) (Conc=100.00) (Conc=100.00) (Conc=100.00) (Conc=100.00) (Conc=100.00) (Conc=100.00) (Conc=100.00) (Conc=100.00) (Conc=100.00) (A.4756 6.59682 3.46 ** ** #* ### ### ### ### ###							
p & m-Xylene 1.41406 1.47559 4.35 (Conc=100.00) o-Xylene 1.43177 1.44755 1.10 Styrene .64562 .65686 1.74 Isopropylbenzene 1.89480 1.93251 1.99 1,1,2,2-Tetrachloroethane .57686 .59682 3.46 ** p-Bromofluorobenzene (SURR) .67535 .69319 2.64 1,2,3-Trichloropropane .16768 .17259 2.93 Bromobenzene .63774 .65675 2.98 n-Propylbenzene .45524 .46106 1.28 2-Chlorotoluene .41764 .44855 7.40 4-Chlorotoluene .45803 .43718 4.55 1,3,5-Trimethylbenzene 1.61365 1.63576 1.37 1,2,4-Trimethylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16	• • •	1.74668	1.74948	.16	*		
0-Xylene	•	1.41406	1.47559	4.35			(Conc=100.00)
Styrene 1.89480 1.93251 1.99 1.1,2,2-Tetrachloroethane 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 3.46 *** 57686 59682 5988 5788 5788 5788 58882 5888	· ·						•
1,1,2,2-Tetrachloroethane p-Bromofluorobenzene (SURR)	•	.64562	.65686	1.74			
p-Bromofluorobenzene (SURR)	Isopropylbenzene	1.89480	1.93251	1.99			
p-Bromofluorobenzene (SURR)	1,1,2,2-Tetrachloroethane	.57686	.59682	3.46		**	
## Bromobenzene		.67535	.69319	2.64			
n-Propylbenzene .45524 .46106 1.28 2-Chlorotoluene .41764 .44855 7.40 4-Chlorotoluene .45803 .43718 4.55 1,3,5-Trimethylbenzene 1.62309 1.61498 .50 tert-Butylbenzene 1.61365 1.63576 1.37 1,2,4-Trimethylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 2.21808 2.22863 .48 1,3-0ichlorobenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.3427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	1,2,3-Trichloropropane	. 16768	.17259	2.93			
2-Chlorotoluene .41764 .44855 7.40 4-Chlorotoluene .45803 .43718 4.55 1,3,5-Trimethylbenzene 1.62309 1.61498 .50 tert-Butylbenzene 1.61365 1.63576 1.37 1,2,4-Trimethylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 2.21808 2.22863 .48 1,3-Dichlorobenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.3427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	Bromobenzene	.63774	.65675	2.98			
4-Chlorotoluene .45803 .43718 4.55 1,3,5-Trimethylbenzene 1.62309 1.61498 .50 tert-Butylbenzene 1.61365 1.63576 1.37 1,2,4-Trimethylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 2.21808 2.22863 .48 1,3-Dichlorobenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.3427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	n-Propylbenzene	.45524	.46106	1.28			
1,3,5-Trimethylbenzene 1.62309 1.61498 .50 tert-Butylbenzene 1.61365 1.63576 1.37 1,2,4-Trimethylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 2.21808 2.22863 .48 1,3-Dichlorobenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.3427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	2-Chlorotoluene	.41764	. 44855	7.40			
tert-Butylbenzene 1.61365 1.63576 1.37 1,2,4-Trimethylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 2.21808 2.22863 .48 1,3-Dichlorobenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.1822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.3427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	4-Chlorotoluene	.45803	.43718	4.55			
1,2,4-Trimethylbenzene 1.56897 1.57045 .09 sec-Butylbenzene 2.21808 2.22863 .48 1,3-Dichlorobenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.3427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	1,3,5-Trimethylbenzene	1.62309	1.61498	.50			
sec-Butylbenzene 2.21808 2.22863 .48 1,3-Dichlorobenzene 1.03711 1.06074 2.28 p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane .13427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	tert-Butylbenzene	1.61365	1.63576	1.37			
1,3-0ichlorobenzene 1.03711 1.06074 2.28 p-1sopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane 1.3427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	1,2,4-Trimethylbenzene	1.56897	1.57045	.09			
p-Isopropyltoluene 1.80992 1.83815 1.56 1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane .13427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	sec-Butylbenzene						
1,4-Dichlorobenzene 1.12622 1.12822 .18 1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane .13427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	1,3-Dichlorobenzene						
1,2-Dichlorobenzene .98178 .98022 .16 n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane .13427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	p-1sopropyltoluene						
n-Butylbenzene 1.84422 1.82916 .82 1,2-Dibromo-3-Chloropropane .13427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	1,4~Dichlorobenzene	_					
1,2-Dibromo-3-Chloropropane .13427 .13950 3.89 1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	•						
1,2,4-Trichlorobenzene .90669 .92061 1.53 Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	·						
Naphthalene 1.12938 1.10858 1.84 Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35							
Hexachlorobutadiene .86158 .85528 .73 1,2,3-Trichlorobenzene .80946 .79040 2.35	1,2,4-Trichlorobenzene						
1,2,3-Trichlorobenzene .80946 .79040 2.35	r				•		
11=1=							

- Response Factor from daily standard file at 50.00 ppb RF

- Average Response Factor from Initial Calibration Form VI RF

*Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/30/95

Time: 15:55

Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 01/24/95

Minimum RF for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

Compound	RF	RF	%Diff	ccc	SPCC	
Dichlorodifluoromethane	3.28484	3.30927	.74			
Chloromethane	1.23549	1.03118	16.54		**	
Vinyl Chloride	1.18628	1.05360	11.19	*		
Bromomethane	1.12701	1.17117	3.92			
Chioroethane		.68840				
Trichlorofluoromethane		4.30491				
Acetone	.54483					
1,1-Dichloroethene		1.36244		*		
Carbon Disulfide		3.32824				
Methylene Chloride		1.34012				
trans-1,2-Dichloroethene		1.47629				
Methyl-tertiary-butylether		2.66306				
1,1-Dichloroethane		2.66450			**	
2-Butanone	.11828					
2,2-Dichloropropane		2.08375				
cis-1,2-Dichloroethene		1.51471		_		
Chloroform		3.41667		*		
1,1-Dichloropropene		2.14361				
1,2-Dichloroethane		2.08329				
1,2-Dichloroethane-d4 (SURR)		1.79184	5.08			
1,1,1-Trichloroethane	.71675	.77330				
Carbon Tetrachloride	.49380	.55411				
Benzene	.86139					
Trichloroethene	.42868					
1,2-Dichloropropane	.39471		7.06	*		
Dibromomethane	.38934					
Bromodichloromethane	.69009					
cis-1,3-Dichloropropene	.47041				(Conc=50.	
trans-1,3-Dichloropropene	.40744				(Conc≃50.	.00)
1,1,2-Trichloroethane	.29985		5.67			
1,3-Dichtoropropane	.51560					
Dibromochloromethane	.63713	.66665	4.63			

- Response Factor from daily standard file at 50.00 ppb RF

RF - Average Response Factor from Initial Calibration Form V1 %Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**) RF

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/30/95 Time: 15:55

Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 01/24/95

Minimum RF for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

Compound	RF	RF	% Diff	ccc	SPCC	
1,2-Dibromoethane	.49176	.50563	2.82			
Bromoform	.47935	.49209	2.66		**	
4-Methyl-2-Pentanone	.10816	.09059	16.24			
Toluene-d8 (SURR)	1.24008	1.20837	2.56			
Toluene	.77901	.77101	1.03	*		
2-Hexanone	. 13407	. 11534	13.97			
Tetrachloroethene	. 71258	.76189	6.92			
Chlorobenzene	1.02796	1.02748	. 05		**	
1,1,1,2-Tetrachloroethane	.58267	.60884	4.49			
Ethylbenzene	1.71442	1.66821	2.70	*		
p & m-Xylene	1.44346	1.44399	.04			(Conc=100.00)
o-Xylene	1.40216	1.37114	2.21			
Styrene	.65131	.64632	.77			
l sopropy i benzene	1.83192	1.86224	1.65			
1,1,2,2-Tetrachioroethane	.58477	.53158	9.09		**	
p-Bromofluorobenzene (SURR)	.64185	.66225	3.18			
1,2,3-Trichloropropane	.16387		2.48			
Bromobenzene	.60052	.62622	4.28			
n-Propylbenzene	.42850	.43434	1.36			
2-Chlorotoluene	.40612	.42219	3.96			
4-Chlorotoluene	.43981		4.88			
1,3,5-Trimethylbenzene	1.59684	1.61754	1.30			
tert-Butylbenzene		1.59076	3.30			
1,2,4-Trimethylbenzene	1.54360		.92			
sec-Butylbenzene		2.13806	.41			
1,3-Dichlorobenzene		1.00173	2.77			
p-Isopropyltoluene	1.74636		.64			
1,4-Dichlorobenzene	1.04693		3.29			
1,2-Dichlorobenzene	.91234	.95769	4.97			
n-Butylbenzene		1.82667	.48			
1,2-Dibromo-3-Chloropropane	. 13164		7.42			
1,2,4-Trichlorobenzene	.84978		5.71			
Naphthalene	1.04596		6.42			
Hexachlorobutadiene	.78724		8.85			
1,2,3-Trichlorobenzene	.75573	.77306	2.29			

- Response Factor from daily standard file at 50.00 ppb

RF - Average Response Factor from Initial Calibration Form VI

%Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/31/95 Time: 10:38

Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 01/24/95

Minimum RF for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

Compound	ŔF	RF	%Diff	ccc	SPCC
Dichlorodifluoromethane	3 28484	3.26401	.63		
Chioromethane		1.05335	14.74		**
Vinyl Chloride		1.06341	10.36	*	
Bromomethane		1.18116	4.80		
Chloroethane	.67266	.69678	3.59		
Trichlorofluoromethane	4.05709	4.15515	2.42		
Acetone	.54483	.52082	4.41		
1,1-Dichloroethene	1.38885	1.36316	1.85	*	
Carbon Disulfide	3.61099	3.24964	10.01		
Methylene Chloride	1.40957	1.33552	5.25		
trans-1,2-Dichloroethene	1.49423	1.48100	.89		
Methyl-tertiary-butylether	2.77508	2.84025	2.35		
1.1-Dichloroethane	2.80844	2.61787	6.79		**
2-Butanone	. 11828	. 10639	10.06		
2,2-Dichloropropane	2.01213	2.11909	5.32		
cis-1,2-Dichloroethene	1.52363	1.50501	1.22		
Chloroform	3.34723	3.34187	. 16	*	
1,1-Dichtoropropene	2.22664	2.12946	4.36		
1,2-Dichloroethane	2.15505	2.11884	1.68		
1,2-Dichloroethane-d4 (SURR)	1.88782	1.83302	2.90		
1,1,1-Trichloroethane	.71675	. 7 3 525	2.58		
Carbon Tetrachloride	.49380	.52102	5.51		
Benzene	.86139	.81091	5. 8 6		
Trichloroethene	.42868	.46479	8.42		
1,2-Dichloropropane	.39471	.37642	4.63	*	
Dibromomethane	. 38934	.40067	2.91		
Bromodichloromethane	.69009	.70142	1.64		
cis-1,3-Dichloropropene	.47041	.47411	.79		
trans-1,3-Dichloropropene	.40744	.42183	3.53		
1,1,2-Trichloroethane	.29985	.29040	3.15		
1,3-Dichloropropane	.51560		2.00		
Dibromochloromethane	.63713	.63488	.35		

RF - Response Factor from daily standard file at 50.00 ppb

RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

2 1 9302

Table 2.8 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/31/95

Time: 10:38

Instrument ID: GCMSD-2(3034A12982) Initial Calibration Date: 01/24/95

Minimum RF for SPCC is 0.30

Maximum % Diff for CCC is 25.0%

Compound	RF	RF	%Diff	CCC	SPCC	
1,2-Dibromoethane	.49176	.50485	2.66			
Bromoform	.47935	.48333	.83		**	
4-Methyl-2-Pentanone	.10816	.10380	4.03			
Toluene-d8 (SURR)	1.24008	1.19932	3.29			
Toluene	.77901	.75518	3.06	*		
2-Hexanone	. 13407	.12971	3.25			
Tetrachloroethene	.71258	.72595	1.88			
Chlorobenzene	1.02796	1.01074	1.68		**	
1,1,1,2-Tetrachloroethane	.58267	.57832	.75			
Ethylbenzene	1.71442	1.66809	2.70	*		
p & m-Xylene	1.44346	1.44396	.03			(Conc=100.00)
o-Xylene	1.40216	1.38607	1.15			
Styrene	.65131	.64868	.40			
Isopropylbenzene	1.83192	1.83271	.04			
1,1,2,2-Tetrachloroethane	.58477	.54251	7.23		**	
p-Bromofluorobenzene (SURR)	.64185	.67239	4.76			
1,2,3-Trichloropropane	. 16387	.16581	1.18			
Bromobenzene	.60052	.61848	2.99			
n-Propylbenzene	.42850	.43548	1.63			
2-Chlorotoluene	.40612	.41388	1.91			
4-Chlorotoluene	.43981	.42890	2.48			
1,3,5-Trimethylbenzene	1.59684	1.59268	.26			
tert-Butylbenzene	1.53990	1.57459	2.25			
1,2,4-Trimethylbenzene	1.54360	1.53225	.74			
sec-Butylbenzene	2.14677	2.13756	.43			
1,3-Dichlorobenzene	.97477	.99352	1.92			
p-Isopropyltoluene	1.74636	1.79593	2.84			
1,4-Dichlorobenzene	1.04693	1.06849	2.06			
1,2-Dichlorobenzene	.91234	.92052	.90			
n-Butylbenzene	1.81789		1.80			
1,2-Dibromo-3-Chloropropane	. 13164	.12764	3.04			
1,2,4-Trichlorobenzene	. 84978	.86661	1.98			
Naphthalene	1.04596		2.63			
Hexach Lorobutadiene		.83425	5.97			
1,2,3-Trichlorobenzene	.75573	.74815	1.00			
•		• • • • • • • • • • • • • • • • • • • •				

RF - Response Factor from daily standard file at 50.00 ppb

RF - Average Response Factor from Initial Calibration Form VI

*Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)

2) 0303

Table 2.9 Correlation of Samples and Calibrations

Samples	Initial Calibration	Continuing Calibration
None	11/22/94	
Lab Blank, 00638, 00635 MS, 00635 MSD	11/22/94	1/13/95
Lab Blank-Water, Lab Blank-Soil, 00640, 00643 00644, 00642 MS, 00642 MSD	11/22/94	1/14/95
Lab Blank-Soil, 00642	11/22/94	1/17/95
Lab Blank, 00636, 00641, 00635, 00639, 00634, 00637	1/18/95	
Lab Blank-Water, Lab Blank-Soil, 00639	1/18/95	1/18/95
None	1/24/95	
Lab Blank, 00914, 00912, 00913, 00914 MS, 00914 MSD	1/24/95	1/30/95
Lab Blank, 00914, 00912, 00913, 00912, 00913	1/24/95	1/31/95

Table 2.10 Results of the TCLP Surrogate Recoveries WA # 0~098 Pier Drum

Sample ID	S1 (DIC)	S2 (TOL)	S3 (BFB)	Total Out	
Lab Blank	98	98	102	0	
Blank Spike	113	100	107	0	
95-01-227-1 MS	105	103	105	0	
95-01-227-1 MSD	115 *	102	106	1	
B 00644	115 *	101	108	1	
B 00643	102	101	107	0	
B 00642	107	101	105	0	

	QC Limits
S1 (DIC) = 1,2-Dichloroethane-d4	76-114
S2 (TOL) = Toluene~d8	88-110
S3 (BFB) = Bromofluorobenzene	86-115

Table 2.11 Results of the TCLP Internal Standard Areas WA # 0-098 Pier Drum

Sample ID	ISTD 1 AREA	ISTD 2 AREA	ISTD 3 AREA	Total Out
50 ppb Std	17237	66998	50342	
95-01-227-1 MS	18092	54895	49536	0
95-01-227-1 MSD	12618	62303	35123	0
Blank Spike	13039	44370	36043	. 0
50 ppb Std	16162	69192	50749	
B 00643	14388	52563	39712	0
B 00642	14158	52247	40860	0
B 00644	14566	52968	40636	0

the acceptable QC range is 50 - 200% of the appropriate standard

ISTD 1 denotes bromochloromethane ISTD 2 denotes 1,4-difluorobenzene

ISTD 3 denotes chlorobenzene-d5

Table 2.12 Results of the MS/MSD Analysis for the TCLP VOA Samples WA # 0-098 Pier Drum

Sample ID: 95-01-227-1

Parameter	Spike Added	Sample Conc.	MS Recov.	~ %	MSD Recov.	%			QC mits
, enditates	(ug/L)	(ug/L)	(ug/L)	Recovery	(ug/L)	Recovery	RPD	RPD	Rec.
1,1-Dichloroethene	500	ND	336.3	67	317.7	64	6	14	61-145
Trichloroethene	500	ND	400.0	80	380.4	76	5	14	71-120
Benzene	500	ND	410.8	82	409.5	82	0	11	76-127
Toluene	500	ND	425.7	8 5	408.5	82	4	13	76-125
Chlorobenzene	500	ND	430.9	8 6	428.5	86	1	13	75-130

Table 2.13 Results of the TCLP Surrogate Recoveries for the Drum Solid Samples

WA # 0-098 Pier Drum

Sample ID	S1 (DIC)	S2 (TOL)	\$3 (BFB)	Total Out
Lab Blank	84	102	96	0
Lab Blank MS	81	100	102	0
00911 MS	79	103	111	0
00911 MSD	81	102	102	0
00911	85	100	102	0
00915	91	103	110	0
00916	90	100	108	0

20	L	mits

S1 (DIC) = 1,2-Dichloroethane-d4	76-114
S2 (TOL) = Toluene-d8	88-110
S3 (BFB) = Bromofluorobenzene	86-115

Table 2.14 Results of the TCLP Internal Standard Areas for the Drum Solid Samples

WA # 0-098 Pier Drum

Sample ID	ISTD 1 AREA	ISTD 2 AREA	ISTD 3 AREA	Total Out
50 ppb Std	70233	427523	341877	
Blank	81241	471191	365543	0
00911	68988	403569	315750	0
00911 MS	74163	442206	380584	0
00911 MSD	74929	434047	360285	0
Blank Spike	73923	426348	342345	0
50 ppb Std	57300	347449	280350	
00911	66428	392236	313525	0
00915	63039	350996	285025	0
00916	62723	347226	279306	0

the acceptable QC range is 50 - 200% of the appropriate standard

ISTD 1 denotes bromochloromethane

ISTD 2 denotes 1,4-diffuorobenzene

ISTD 3 denotes chlorobenzene-d5

Table 2.15 Results of the Blank Spike Analysis for the TCLP Analysis of the Drum Solid Samples WA # 0-098 Pier Drum

Compound	Sample Conc (ug/l)	Spike Conc (ug/l)	Recovered Conc (ug/l)	% Recovery (ug/l)	QC Limits	2 - 2	
1,1-Dichloroethene	24	20	48	120	61-145		
Benzene	25	20	50	125 *	71-120		
Trichloroethene	21	20	42	105	76-127		
Toluene	24	20	48	120	76-125		
Chlorobenzene	24	20	48	120	75-130		

Table 2.16 Results of the MS/MSD Analysis for the TCLP VOA Analysis of the Drum Solid Samples WA # 0 \sim 098 Pier Drum

Sample ID: 00911

Parameter	Spike Added	Sample Conc.	MS Recov.	%	MSD Recov.	%			QC mits
	(ug/L)	(ug/L)	(u g /L)	Recovery	(ug/L)	Recovery	RPD	RPD	Rec.
1,1-Dichloroethene	100000	ND	113360	113	112040	112	1	14	61-145
Trichloroethene	100000	ND	96780	97	97140	97	0	14	71-120
Benzene	100000	ND	114140	114	115080	115	1	11	76~127
Toluene	100000	3140	110880	108	111220	108	0	13	76-125
Chlorobenzene	100000	ND	109100	109	113700	114	4	13	75-130

2 1 0101

QA/QC FOR BNA

Before extraction, each sample was spiked with a six component mixture of CLP surrogate standards consisting of nitrobenzene-d₅, 2-fluorobiphenyl, terphenyl-d₁₄, phenol-d₅, 2-fluorophenol, and 2,4,6-tribromophenol. The surrogate percent recoveries for the water samples, listed in Table 2.17, ranged from 27 to 125. Fifty-seven values out of sixty were within the acceptable QC limits.

The internal standard areas (for 1,4-dichlorobenzene- d_4 , naphthalene- d_8 , acenaphthene- d_{10} , phenanthrene- d_{10} , chrysene- d_{12} , perylene- d_{12}) for the water samples are also listed in Table 2.17. All ninety-six areas are within the acceptable QC limits.

Samples CD 00640 and BE 00527 were chosen for the matrix spike/matrix spike duplicate (MS/MSD) analyses for the water samples. The percent recoveries, ranging from 37 to 84, are listed in Table 2.18. Twenty-one values out of twenty-two were within the acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.18, ranged from 2 to 18 and all 11 values were within the acceptable QC limits.

The surrogate percent recoveries for the soil samples, listed in Table 2.19, ranged from 70 to 109. All thirty-six values were within the acceptable QC limits.

The internal standard areas for the soil samples are also listed in Table 2.19. All thirty-six areas are within the acceptable QC limits.

Sample A 00644 was chosen for the MS/MSD analyses for the soil samples. The percent recoveries, ranging from 69 to 108, are listed in Table 2.20. Eighteen out of twenty-two values were within the acceptable QC limits. The RPDs, also listed in Table 2.20, ranged from 1 to 20 and all eleven values were within the acceptable QC limits.

The surrogate percent recoveries for the TCLP samples, listed in Table 2.21, ranged from 40 to 117. All thirty values were within the acceptable QC limits.

The internal standard areas for the TCLP samples are also listed in Table 2.21. All thirty areas are within the acceptable QC limits.

Sample BE 00527 was chosen for the MS/MSD analyses for the TCLP samples. The percent recoveries, ranging from 61 to 104, are listed in Table 2.22. Nineteen out of twenty-two values were within the acceptable QC limits. The RPDs, also listed in Table 2.22, ranged from 0 (zero) to 29 and ten out of eleven values were within the acceptable QC limits.

The surrogate percent recoveries for the drum liquid samples, listed in Table 2.23, ranged from 33 to 168. Thirty out of thirty-eight values were within the acceptable QC limits. Twenty-two other values were from diluted samples and the percent recoveries were not calculated.

The internal standard areas for the drum liquid samples are also listed in Table 2.23. Fifty-six out of sixty areas are within the acceptable QC limits.

Water blank WBLK021495 was chosen for the MS/MSD analyses for the drum liquid samples. The percent recoveries, ranging from 34 to 82, are listed in Table 2.24. All twenty-two values were within the acceptable QC limits. The RPDs, also listed in Table 2.24, ranged from 0 (zero) to 14 and all eleven values were within the acceptable QC limits.

The results of the initial calibration are listed in Table 2.25.

The results of the continuing calibration are listed in Table 2.26.

The samples and calibrations are correlated in Table 2.27.

The surrogate percent recoveries for the TCLP extracts of the drum solid samples, listed in Table 2.28, ranged from 11 to 111. Fifty out of fifty-one values were within the acceptable QC limits. Three other values were not calculated because of matrix interference in the sample. The value that exceeded the acceptable QC limit was also affected by matrix interference.

The internal standard areas for the TCLP extracts of the drum solid samples are listed in Table 2.29. Fifty-three out of fifty-four areas are within the acceptable QC limits.

The results of the blank spike analysis are listed in Table 2.30. The percent recoveries ranged from 42 to 150 and all eleven values were within the acceptable QC limits.

Sample 00915 was chosen for the MS/MSD analyses for the TCLP extracts of the drum solid samples. The percent recoveries, ranging from 12 to 95, are listed in Table 2.31. Twenty-one out of twenty-two values were within the acceptable QC limits. The RPDs, also listed in Table 2.31, ranged from 2 to 59 and seven out of eleven values were within the acceptable QC limits.

2 / 0403

Table 2.17 Results of the Internal Standard Areas and Surrogate Recoveries for the Water Samples WA # 0-098 Pier Drum

CAL CHECK 50 PPM BNA	>PD007	39875 250541	174 38 4 2501 8 5	129444 296069	NA NA	NA NA	NA NA
WBLK011895	>PD010	31905 198702	142809 240 96 7	102084 214443	27* 59	30 * 70	46 76
CD00636	>PD011	30721 20 78 42	133957 258949	95093 242135	75 71	85 79	62 94
CD00638	>PD012	33455 205400	145356 284653	124516 239478	94 84	88 102	69 125*
CD00641	>PD013	33306 230724	151794 294088	114365 254581	84 79	76 94	62 101
CD00635	>PD014	38167 238915	171288 292358	125508 274031	88 67	93 80	64 103
CD00639	>PD015	33135 228271	151576 255086	117413 248236	77 70	71 92	69 96
CD00640	>PD016	33189 202208	125163 228261	91698 234704	59 70	58 84	58 81
CD00640 MS	>PD017	32712 204117	138868 271469	107572 2 3868 9	57 73	56 8 5	53 90
CD00640 MSD	>PD018	32325 204579	135612 282333	107602 258520	49 77	48 80	46 88
CD00634	>PD019	34238 235951	153494 292426	116291 264288	91 90	90 94	68 97

		SI	JRROGATE LIMITS	WATER
S1	(NBZ)	=	Nitrobenzene-d5	(35-114)
S2	(FBP)	=	2-Fluorobiphenyl	(43-116)
S 3	(TPH)	=	Terphenyl-d14	(35-114)
\$4	(PHL)	=	Phenol-d5	(10-110)
S 5	(2FP)	=	2-fluorophenol	(21-110)
S6	(TBP)	=	2,4,6-Tribromophenol	(10-123)

Table 2.18 Results of the MS/MSD Analysis for the Water Samples WA # 0-098 Pier Drum

Matrix Spike - EPA Sample No.: CD00640

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (Ug/L)	MS CONCENTRATION (Ug/L)	MS % REC #	QC LIMITS REC.
		=========		======	
Phenol	156.25	0.00	111.96		12- 89
2-Chlorophenol	156.25	0.00	115.54	73	27-12
1,4-Dichlorobenzene	78.13	0.00	32.21	41	36- 9
N-Nitroso-di-n-prop.(1)	78.13	0.00	40.69	52	41-11
1,2,4-Trichlorobenzene	78.13	0.00	33.85	43	39- 9
4-Chloro-3-methylphenol	156.25	0.00	120.71	77	23- 9
Acenaphthene	78.13	0.00	46.43	59	46-11
4-Nitrophenol	156.25	0.00	132.21	84 *	10- 8
2,4-Dinitrotoluene	78.13	0.00	49.25	63	24- 9
Pentachlorophenol	156.25	0.00	126.99	81	9-10
Pyrene	78.13	0.00	52.05	66	26-12

ADDED (ug/L) ======== 156.25	CONCENTRATION (ug/L)	% REC #	% RPD #	QC L	REC.
75 25	=======================================	=====!			:
	107.20	68	-4	42	12- 89
			2		27-123
		37	10		36- 97
78.13	37.15	47	10	38	41-116
78.13	34.41	44	2	28	39- 98
156.25	131.69	84	8	42	23- 97
78.13	44.59	57	3	31	46-118
156.25	123.95	79	6	50	10- 80
78.13	47.13	60	4	38	24- 96
156.25	121.33	77	5	50	9-103
78.13	43.07	55	18	31	26-127
	78.13 78.13 78.13 156.25 78.13 156.25 78.13	78.13 29.13 78.13 37.15 78.13 34.41 156.25 131.69 78.13 44.59 156.25 123.95 78.13 47.13 156.25 121.33	78.13 29.13 37 78.13 37.15 47 78.13 34.41 44 156.25 131.69 84 78.13 44.59 57 156.25 123.95 79 78.13 47.13 60 156.25 121.33 77	78.13 29.13 37 10 78.13 37.15 47 10 78.13 34.41 44 2 156.25 131.69 84 8 78.13 44.59 57 3 156.25 123.95 79 6 78.13 47.13 60 4 156.25 121.33 77 5	78.13 29.13 37 10 28 78.13 37.15 47 10 38 78.13 34.41 44 2 28 156.25 131.69 84 8 42 78.13 44.59 57 3 31 156.25 123.95 79 6 50 78.13 47.13 60 4 38 156.25 121.33 77 5 50

⁽¹⁾ N-Nitroso-di-n-propylamine

7 . 3.03

Table 2.19 Results of the Internal Standard Areas and Surrogate Recoveries for the Soil Samples WA # 0-098 Pier Drum

			Inter	nal Standard	is	Sur	rogates	i
			1	2	3	NBZ	FBP	TPH
		Data	4	5	6	PHL	2FP	TBP
	Sample #	file	area	area	area	*	*	*
CAL CHECK	50 PPM BNA	>PD021	38434	163272	120938	NA	NA	NA
			231504	234487	248960	NA	NA	NA
	SBLK011395	>PD024	36298	150688	115900	92	96	65
			211649	278112	220347	87	87	99
	A00642	>PD025	34691	152529	102508	80	102	77
		, , , ,	212295	243231	210899	79	90	104
	A00643	>PD026	32854	138441	103061	96	101	76
			201910	229956	207427	98	106	108
	A00644	>PD027	30669	142742	106559	77	84	70
		, , , ,	196969	241902	203685	85	79	97
	A00644 MS	>PD028	31675	135866	97579	84	99	80
			189961	228494	200295	89	95	109
	A00644 MSD	>PD029	26902	120589	99910	98	88	84
			191646	224721	192040	92	109	99

	SOIL			
\$1	(NBZ)	=	Nitrobenzene-d5	(23-120)
S2	(FBP)	=	2-Fluorobiphenyl	(30-115)
\$3	(TPH)	=	Terphenyl-d14	(18 137)
\$4	(PHL)	=	Phenot-d5	(24-113)
S 5	(2FP)	=	2-Fluorophenol	(25-121)
56	(TBP)	=	2,4,6-Tribromophenol	(19-122)

Table 2.20 Results of the MS/MSD Analysis for the Soil Samples WA # 0-098 Pier Drum

Matrix Spike - EPA Sample No.: A00644

Level:(low/med) LOW

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (Ug/Kg)	MS % REC #	QC LIMITS REC.
Phenol	4065.04	0.00	4031.18	99 *	26- 90
2-Chlorophenol	4065.04	0.00	3646.30	89	25-102
1,4-Dichlorobenzene	2032.52	0.00	1573.13	77	28-104
N-Nitroso-di-n-prop.(1)	2032.52	0.00	1562.72	76	41-126
1,2,4-Trichlorobenzene_	2032.52	0.00	1734.39	85	38-107
4-Chioro-3-methylphenol	4065.04	0.00	3996. 10	98	26-103
Acenaphthene	2032.52	0.00	1900.73	93	31-137
4-Nitrophenol	4065.04	0.00	4204.63	103	11-114
2,4-Dinitrotoluene	2032.52	0.00	2018.82	99 *	28- 89
Pentachlorophenol	4065.04	0.00	2808.33	69	17-109
Pyrene	2032.52	0.00	1859.76	91	35-142

!	SPIKE ADDED	MSD CONCENTRATION	MSD %	*	QC L	IMITS
COMPOUND	(ug/Kg)	(ug/Kg)	REC #	RPD #	RPD	REC.
	======		======	=====	=====	=====
Phenoi	4065.04	4403.00	108 *	11	35	26- 90
2-Chlorophenol	4065.04	4025.73	99	10	50	25-102
1,4-Dichlorobenzene	2032.52	1813.58	89	14	27	28-104
N-Nitroso-di-n-prop.(1)	2032.52	1956.38	96	23	38	41-126
1,2,4-Trichlorobenzene	2032.52	1900.85	93	8	23	38-107
4-Chloro-3-methylphenol	4065.04	4116.00	101	5	33	26-103
Acenaphthene	2032.52	1874.71	92	1	19	31-137
4-Nitrophenol	4065.04	3982.72	97	6	50	11-114
2,4-Dinitrotoluene	2032.52	1871.91	92 *	7	47	28- 89
Pentachlorophenol	4065.04	3473.37	85	20	47	17-109
Pyrene	2032.52	1780.00	87	4	36	35-142

⁽¹⁾ N-Nitroso-di-n-propylamine

(· · ·)

Table 2.21 Results of the Internal Standard Areas and Surrogate Recoveries for the TCLP Samples WA # 0-098 Pier Drum

			Inter	nal Standard	is	Sur	rogates	•
	Sample #	Data File	1 4	2 5	3 6	NBZ PHL %	FBP 2FP %	TPH TBP
	Sample #		area	area	area			
CAL CHECK	50 PPM BNA	>SM005	41021	170697	120567	NA	NA	NA
			241360	258471	300823	NA	NA	NA
	WBLK011995	>PD001	38169	144186	115204	81	75	66
			239769	298718	281582	40	48	82
	воо525	>PD002	46924	212640	159659	80	69	71
			290233	364226	351123	55	68	99
	B00526	>PD003	38582	175233	111491	77	81	74
			218230	280286	276763	46	69	98
	BE00527	>PD004	41094	165496	129866	75	66	65
			228236	291819	289160	54	59	91
CAL CHECK	50 PPM BNA	>PD007	39875	174384	129444		NA	 NA
			250541	250185	296069	NA	NA	NA
	BE00527 MS	>PD008	41269	199129	146879	93	93	67
			279101	389251	339551	94	100	117
	BE00527 MSD	>PD009	46081	192055	144902	89	82	72
			254658	328287	343633	77	87	115

		SI	JRROGATE LIMITS	WATER
S1	(NBZ)	=	Nitrobenzene-d5	(35-114)
S2	(FBP)	=	2-Fluorobiphenyl	(43-116)
S 3	(TPH)	=	Terphenyl-d14	(35-114)
54	(PHL)	=	Phenoi -d5	(10-110)
S 5	(2FP)	=	2-Fluorophenol	(21-110)
S6	(TBP)	Ξ	2,4,6-Tribromophenol	(10-123)

2 ... 94.23

Table 2.22 Results of the MS/MSD Analysis for the TCLP Samples WA # 0-098 Pier Drum

Matrix Spike - EPA Sample No.: BE00527

COMPOUND	SPIKE	SAMPLE	MS	MS	QC
	ADDED	CONCENTRATION	CONCENTRATION	%	LIMITS
	(ug/L)	(ug/L)	(ug/L)	REC #	REC.
Phenol 2-Chlorophenol 1,4-Dichlorobenzene N-Nitroso-di-n-prop.(1) 1,2,4-Trichlorobenzene 4-Chloro-3-methylphenol Acenaphthene 4-Nitrophenol 2,4-Dinitrotoluene Pentachlorophenol Pyrene	163.93 163.93 81.97 81.97 81.97 163.93 81.97 163.93 81.97	0.00 0.00 0.00 0.00 0.00 0.00 0.00	67.38 79.75 62.72 143.73 72.39 140.03 78.62 164.66	82 97 76 87 88 85 *	12- 89 27-123 36- 97 41-116 39- 98 23- 97 46-118 10- 80 24- 96 9-103 26-127

		¦ MSD ¦	MSD		i	
	ADDED	CONCENTRATION	*	*	QC L	IMITS
COMPOUND	(ug/L)	(ug/L)	REC #	RPD #	RPD	REC.
*****************	=======	=============	=====	=====	=====	=====
Phenol	163.93	119.72	73	17	42	12- 89
2-Chlorophenol	163.93	135.11	82	20	40	27-123
1,4-Dichlorobenzene	81.97	50.14	61	29 *	28	36- 97
N-Nitroso-di-n-prop.(1)	81.97	63.38	77	22	38	41-116
1,2,4-Trichlorobenzene	81.97	52.04	63	18	28	39- 98
4-Chloro-3-methylphenol	163.93	132.14	80	8	42	23- 97
Acenaphthene	81.97	75.48	92	4	31	46-118
4-Nitrophenol	163.93	147.32	89 *	4	50	10- 80
2,4-Dinitrotoluene	81.97	78.42	95	0	38	24- 96
Pentachlorophenol	163.93	170.63	104 *	3	50	9-103
Pyrene	81.97	67.12	81	9	31	26-127
-	i	i			i	i

⁽¹⁾ N-Nitroso-di-n-propylamine

2 1 9409

Table 2.23 Results of the Internal Standard Areas and Surrogate Recoveries for the Drum Liquid Samples WA # 0-098 Pier Drum

			Internal Standards			Surr	ogates	
			1	2	3	NBZ	FBP	TPH
		Data	4	5	6	PHL	2FP	TBP
	Sample #	File	area	8168	area	×	×	X
CAL CHECK	50 PPM BNA	>PD050	27207	123057	99453	NA.	MA	
CAL CHECK	DU PPM BMA	77000	27207 167961	186638	8865 2 180600	NA NA	NA NA	NA NA
								710
	WBLK021495	>PD051	34901	152328	112436	64	6 6	69
			203720	257540	219488	33		92
	WBLK021495	>PD052	40548	177572	122521	69	79	65
	MS		246256	303794	274396	35	54	91
			/ OFFE	470000	407000			
	WBLK021495	>10055	42555	178299	127220	74 77	79 EE	68 107
	MSD		226030	289916	248586	34	55	107
	00912	>PD054	58615*A	223587	170681	D	117*	91
	007.12		352873*A	364098	332488	Ď	D	Ď
	00913	>PD055	50428	188345	132366	D	D	D .
			244243	249876	215022	D	D	D 🤄
	00914	>PD056	63403*A	143687	167081	168*	-	67
			311429	379043*A	293695	41	63	128*
CAL CHECK	50 PPM BNA	>PD057	36440	144315	101172	NA	NA	NA
			196978	233759	215437	NA	NA	NA
	00912 10X	>PD058	59615	252485	161749	D	D	D
			307077	398538	336595	D	D	D
		>PD059	55289	249859	147017		D	D
	00913 10X	770039	317671	388136	162817 275650	Ð D	D	D
					273030	· · · · · · · · · · · · · · · · · · ·		
	00914 10X	>PD060	50485	197498	139658	115*	106	88
			252995	342917	281060	55	80	161*

CAL CHECK	50 PPM BNA	>PD033	52698 700038	217320	145371	NA	NA	NA
			300028	275807	341122	NA.	NA	NA
	00913 (BOT		45989	201163	149527	113	106	79
	00715 (501		272996	323814	284291	111*	117*	123*
			2. E E					
S	URROGATE LIM	ITS	SOIL					
		JF	437 430:					
	≈ Nitrobenze		(23-120)					
	≈ 2-fluorobi = Terphenyl-		(30-115) (18-137)					
	≈ Terphenyl- ≈ Phenol-d5	W 14	(24-113)					
	= 2-fluoroph	enol	(25-121)					
	= 2,4,6-Trib							
	, ,	,	••					

2 2 3410

Table 2.24 Results of the MS/MSD Analysis for the Drum Liquid Samples WA # 0-098 Pier Drum

Matrix Spike - EPA Sample No.: WBLK021495

	SPIKE ADDED	SAMPLE CONCENTRATION	MS CONCENTRATION	MS %	QC LIMITS
COMPOUND	(ug/L)	(ug/L)	(ug/L)	REC #	REC.
	========		=========	IXSERE	22222
Phenol	100.00	0.00	35.33	35	12- 89
2-Chlorophenol	100.00	0.00	69.64	69	27-123
1,4-Dichlorobenzene	50.00	0.00	28.44	56	36- 97
N-Nitroso-di-n-prop.(1)	50.00	0.00	36.47	72	41-116
1,2,4-Trichlorobenzene	50.00	0.00	28.00	56	39- 98
4-Chloro-3-methylphenol	100.00	0.00	77.21	77	23- 97
Acenaph thene	50.00	0.00	38.73	77	46-118
4-Nitrophenol	100.00	0.00	34.33	34	10- 80
2,4-Dinitrotoluene	50.00	0.00	39.86	79	24- 96
Pentachlorophenol	100.00	0.00	71.41	71	9-103
Pyrene	50.00	0.00	35.61	71	26-127

	SPIKE ADDED	MSD CONCENTRATION	MSD %	x	00.1	MITS
COMPOUND	(ug/L)	(ug/L)	REC #			REC.
	=======	=======================================	=====	=====	=====	=====
Phenol	100.00	34.70	34	2	42	12- 89
2-Chiorophenol	100.00	71.36	71	2	40	27-123
1.4-Dichlorobenzene	50.00	27.87	55	1	28	36- 97
N-Nitroso-di-n-prop.(1)	50.00	34.29	68	5	38	41-116
1,2,4-Trichlorobenzene	50.00	29.83	59	5	28	39- 9
4-Chloro-3-methylphenol	100.00	77.82	77	0	42	23- 9
Acenaphthene	50.00	41.05	82	6	31	46-11
4-Nitrophenol	100.00	35.24	35	2	50	10- 8
2.4-Dinitrotoluene	50.00	38.89	77	2	38	24- 90
Pentachlorophenol	100.00	82.57	82	14	50	9-10
Pyrene	50.00	37.09	74	۱ ۷	31	26-12

⁽¹⁾ N-Nitroso-di-n-propylamine

Table 2.25 Results of the Initial Calibration WA # 0-098 Pier Drum

Instrument ID: 888632 Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % RSD for CCC is 30%

Laboratory ID: Compound	* >GL090 RF 20.00	>GL089 RF 50.00	>GL088 RF 80.00	>GL087 RF 120.00	>GL086 RF 160.00	RRT	RF	% RSD	CCC	SPCC
2-Fluorophenol	1.13310	1.22413	1.11533	1.02194	1.25164	.681	1.14925	7.988		
Phenol-d5			_	1.65337			1.76675	5.046		
Phenol				1.95778			2.13426	5.010	*	
bis(-2-Chloroethyl)Ether				1.44631			1.61547	10.295		
2-Chlorophenol				1.24925			1.49216	10.534		
1,3-Dichlorobenzene				1.35242		-	1.51267	10.932		
1,4-Dichlorobenzene				1.36267			1.55768	11.722	*	
Benzyl alcohol				1.45491			1.58553	7.471		
1,2-Dichlorobenzene				1.16091			1.39231	17,140		
2-Methylphenol	1.38898	1.46332	1,42301	1.19653	1.25385	1.096	1.34514	8.507		
bis(2-Chloroisopropyl)ether	3.99327	3.90325	4.01262	3.40084	3.57889	1.100	3.77777	7.237		
4-Methylphenol	1.66697	1.41573	1,24256	1.02896	1.25535	1.141	1.32191	17.918		
N-Nitroso-Di-n-propylamine	2.30631	1.97411	1.85097	1.82601	2.01124	1.143	1.99373	9.612		**
Hexachloroethane	.56066	.56020	.47790	.42972	.47105	1.140	.49991	11.651		
Nitrobenzene-d5	.44382	.44849	.41265	.42927	.41924	.853	.43070	3.574		
Nitrobenzene	.51451	.50006	.46011	.45953	.43566	.857	.47397	6.830		
Isophorone	1.04231	.99043	.92646	.98080	.99973	.909	.98795	4.214		
2-Nitrophenol	.25557	. 28605	.23798	.24043	.22613	.925	.24923	9.265	*	
2,4-Dimethylphenol	.33113	.36608	.32878	.33612	.30239	.944	.33290	6.821		
bis(2-Chloroethoxy)methane	.59820	.57447	.51769	.50943	.52076	.964	.54411	7.289		
2,4-Dichlorophenol	.36826	.35234	.33708	.34390	.30268	.977	.34085	7.133	*	
1,2,4-Trichlorobenzene	.36276	.39848	.33845	.31560	.30452	.993	.34396	10.990		
Naphthalene	1.25299	1.14715	1.02734	.91575	.95318	1.005	1.05928	13.193		
4-Chloroaniline	.56557	.57445	.50288	.52249	.45405	1.028	.52389	9.358		
Hexach Lorobutadiene	.21860	.22318	. 18806	. 19862	. 16995	1.048	. 19968	11.001	*	
4-Chloro-3-methylphenol	.47239	.45019	.42508	.39662	.38289	1.141	.42543	8.683	*	
2-Methylnaphthalene	.74715	. 75329	.61199	. 59383	.58603	1.157	.65846	12.806		•
Hexachlorocyclopentadiene	.34624	.32617	.36041	.32367	.36697	.869	.34469	5.675		**
2,4,6-Trichlorophenol	.45783	.47252	.44975	.36503	.42045	.884	.43312	9.820	*	
2,4,5-Trichlorophenol	.51589	.52531	.52369	.38522	.42932	.889	.47589	13.584		

RF - Response Factor (Subscript is amount in ng/ul)

RRT - Average Relative Retention Time (RT Std/RT Istd)

- Average Response Factor - Percent Relative Standard Deviation %RSD -- Calibration Check Compounds (*) CCC

SPCC - System Performance Check Compounds (**)

Table 2.25 (Cont) Results of the Initial Calibration WA # 0-098 Pier Drum

Instrument ID: 888632 Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % RSD for CCC is 30%

Laboratory ID:	>GL090 RF	>GL089 RF	>GL088 RF	>GL087 RF	>GL086 RF					
Compound	20.00	50.00	80.00	120.00	160.00	RRT	RF	% RSD	CCC	SPCC
2-Chloronaphthalene	1.34685	1.29151	1.15291	.95044	1.02261	.909	1.15286	14.695	•••	
2-Fluorobiphenyl	1.39267	1.24336	1.25919	.94988	1.01710	.897	1.17244	15.653		
2-Nitroaniline	.60778	.57489	.58188	.50323	.53797	.936	.56115	7.285		
Dimethylphthalate				1.37395			1.59849	15.157		
Acenaphthylene	2.51035	2.12355	2.06257	1.47625	1.51203	.975	1.93695	22.684		-
3-Nitroaniline	.81753	.81116	.79192	.64567	.66796	1.003	.74685	11,127		
Acenaphthene		1.41583	1.15309	.97746	.94646	1.006	1.19323	20.402	*	
2,4-Dinitrophenol	. 25361	.30462	.32098	.31292	.33168	1.018	.30476	9.940		**
4-Nitrophenol	.51513	.50113	.50027	.42020	.44969	1.034	.47728	8.478		**
Dibenzofuran				1.51699				15.413		
2,6-Dinitrotoluene	.43148	.42552	.45502	.36978	.39788	. 983	.41593	7.899		
2,4-Dinitrotoluene	.67089	.65597	.68474	.54380	.63087	1.043	.63725	8.776		
Diethylphthalate				1.01751			1.48656	31.079		
4-Chlorophenyl-phenylether	.91730	.86277	. 76036	.65770	.60591	1.089	.76081	17.327		
fluorene		1.40943		.97477	.96797		1.22830	20.929		
4-Nitroaniline	.59544	.65004	.61573	.54895	.57137	1.102	.59631	6.565		
4,6-Dinitro-2-methylphenol	.20371	. 20695	.22258	.20540	.19005	.900	.20574	5.624		
N-Nitrosodiphenylamine	.70111	. 55363	.53611	.41026	.33753	.903	.50773	27.640	*	
2,4,6-Tribromophenol	. 20633	.17349	.17572	. 15391	. 15263	.916	.17241	12.627		
4-Bromophenyl-phenylether	.32866	.25835	. 25454	. 24703	.21657	.946	.26103	15.791		
Hexachlorobenzene	.40076	. 33396	.31531	.30145	.28440	.963	.32717	13.748		
Pentachlorophenol	.27075	. 25754	. 22043	.22780	.21471	.987	.23825	10.300	*	
Phenanthrene		1.13432		.95399	.92722		1.10818	18.749		
Anthracene		1.27836		.97674	.93599		1.16825	20.227		
Carbazole				1.00621			1.17462	15.431		
Di-n-butylphthalate				1.38343			1.69660	23.914		
fluoranthene				1.16479			1.39808	18.202	*	
Pyrene				1.49885			1.54199	5.285		
Terphenyl-d14				1.11861			1.16385	7.436		
Butylbenzylphthalate	.91649	.82759	.90642	.90127	.97141	.954	.90464	5.677		
3,31-Dichlorobenzidine	.57520	.56042	.57350	.53587		1.000	.56015	2.842		
Benzo(a)anthracene				1.49664			1.57436	7.550		
Bis(2-Ethylhexyl)phthalate				1.12011			1.19680	5.019		
Chrysene				1.26285			1.32497	5.111		
Di-n-octylphthalate					1.56698		1.78652	12.616	*	
Benzo(b)fluoranthene					1.35090		1.49577	8.047		
Benzo(k)fluoranthene		1.15037		.95476			1.07387	24.611	_	
Benzo(a)pyrene				1.25669			1.29324	4.482	*	
Indeno(1,2,3-cd)pyrene					1.33280		1.38492	7.759		
Dibenzo(a,h)anthracene	1.06802	.95635	.84996	.87604	.92602	1.132	.93528	9.094		
Benzo(g,h,i)perylene	1.43203	1.23439	1.10012	1.15408	1.16585	1.10/	1.23341	9.624		

RF - Response Factor (Subscript is amount in ng/ul)

RRT - Average Relative Retention Time (RT Std/RT Istd)

- Average Response Factor

%RSD - Percent Relative Standard Deviation
 CCC - Calibration Check Compounds (*)
 SPCC - System Performance Check Compounds (**)

Table 2.26 Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/27/95 Time: 10:42

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	X Diff	CCC	SPCC
2-Fluorophenol	1.14925	1.08407	5.67		
Phenol-d5		1.65681	6.22		
Phenoi		1.93733	9.23	*	
bis(-2-Chloroethyl)Ether		1.75268	8.49		
2-Chlorophenol		1.54303	3.41		
1,3-Dichlorobenzene		1.50633	.42		
1,4-Dichlorobenzene	1.55768	1.63957	5.26	*	
Benzyl alcohol	1.58553	1.59575	.64		
1,2-Dichlorobenzene	1.39231	1.44697	3.93		
2-Methylphenol	1.34514	1.36088	1.17		
bis(2-Chloroisopropyl)ether	3.77777	3.82251	1.18		
4-Methylphenol	1.32191	1.61184	21.93		
N-Nitroso-Di-n-propylamine		1.91086	4.16		**
Hexachloroethane	.49991		18.61		
Nitrobenzene-d5	.43070		6.65		
Nitrobenzene	.47397		3.53		
Isophorone	.98795		3.96		
2-Nitrophenol	.24923		9.57	*	
2,4-Dimethylphenol	.33290		1.28		
bis(2-Chloroethoxy)methane	.54411		14.11		
2,4-Dichlorophenol	.34085		7.31	*	
1,2,4-Trichlorobenzene	.34396		18.52		
Naphthalene		1.05871	.05		
4-Chloroaniline	.52389				
Hexach Lorobutadiene		.27619	38.32	*	
4-Chloro-3-methylphenol	.42543		_	*	
2-Methylnaphthalene	.65846				
Hexachlorocyclopentadiene	.34469		20.28		**
2,4,6-Trichlorophenol	.43312		18.71	•	
2,4,5-Trichtorophenol	.47589		19.37		
2-Chloronaphthalene		1.24447	7.95		
2-Fluorobiphenyl	1.17244	1.29747	10.66		

- Response Factor from daily standard file at 50.00 ng/ul

RF - Average Response Factor from Initial Calibration Form VI

*ZDiff - % Difference from original average or curve
CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/27/95

Time: 10:42 Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	%Diff	CCC	SPCC
2-Nitroaniline	.56115	.63065	12.38		
Dimethylphthalate		1.79658	12.39		
Acenaphthylene		2.00684	3.61		
3-Nitroaniline	.74685		8.96		
Acenaphthene		1.28618	7.79	*	
2,4-Dinitrophenol		.29768			**
4-Nitrophenol	.47728		3.89		**
Dibenzofuran	1.84025		2.16		
2.6-Dinitrotoluene	.41593		8.68		
2,4-Dinitrotoluene	.63725	.62790	1.47		
Diethylphthalate		1.64452	10.63		
4-Chlorophenyl-phenylether	.76081	.77358	1.68		
Fluorene	1.22830		7.03		
4-Nitroaniline	.59631	.60980	2.26		
4,6-Dinitro-2-methylphenol	.20574	.22411	8.93		
N-Nitrosodiphenylamine	.50773	.59090	16.38	*	
2,4,6-Tribromophenol	.17241	.20206	17.19		
4-Bromophenyl-phenylether	.26103	.26870	2.94		
Hexachlorobenzene	.32717	.41524	26.92		
Pentachlorophenol	.23825	.26073	9.43	*	
Phenanthrene	1.10818	1.19842	8.14		
Anthracene	1.16825	1.26409	8.20		
Carbazole	1.17462	1.15933	1.30		
Di-n-butylphthalate	1.69660	1.63882	3.41		
Fluoranthene	1.39808	1.50420	7.59	*	
Pyrene	1.54199	1.48470	3.72		
Terphenyl-d14	1.16385	1.14176	1.90		
Butylbenzylphthalate	.90464	.78351	13.39		
3,3'-Dichlorobenzidine	.56015	.48883	12.73		
Benzo(a)anthracene	1.57436	1.45835	7.37		
Bis(2-Ethylhexyl)phthalate	1.19680	1.11442	6.88		
Chrysene	1.32497	1.25968	4.93		
Di-n-octylphthalate	1.78652	1.74477	2.34	*	
Benzo(b)fluoranthene	1.49577	1.57277	5.15		
Benzo(k)fluoranthene		1.18389	10.24		
Benzo(a)pyrene	1.29324		4.20	*	
Indeno(1,2,3-cd)pyrene	1.38492		8.58		
Dibenzo(a,h)anthracene	.93528		2.36		
Benzo(g,h,i)perylene	1.23341	1.34492	9.04		
• • • • • • • • • • • • • • • • • • • •					

- Response Factor from daily standard file at 50.00 ng/ul

RF - Average Response Factor from Initial Calibration Form VI

*Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/30/95

Time: 10:15

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	XD iff	CCC	SPCC
2-Fluorophenol	1.14925	1.05456	8.24		
Phenol-d5		1.74913	1.00		
Phenol	2.13426	1.89127	11.39	*	
bis(-2-Chloroethyl)Ether	1.61547	1.74696	8.14		
2-Chlorophenol	1.49216	1.44837	2.93		
1,3-Dichlorobenzene	1.51267	1.44315	4.60		
1,4-Dichlorobenzene	1.55768	1.72617	10.82	*	
Benzyl alcohol	1.58553	1.56684	1.18		
1,2-Dichtorobenzene	1.39231	1.43396	2.99		
2-Methylphenol	1.34514	1.41859	5.46		
bis(2-Chloroisopropyl)ether	3.77777	3.96260	4.89		
4-Methylphenol	1.32191	1.45029	9.71		
N-Nitroso-Di-n-propylamine	1.99373	2.03123	1.88		**
Hexachloroethane	.49991		20.85		
Nitrobenzene-d5	.43070	. 45598	5.87		
Nitrobenzene	.47397	.48215	1.72		
lsophorone	.98795	.99758	.98		
2-Nitrophenol	. 24923	. 22927	8.01	*	
2,4-Dimethylphenol	.33290	.31393	5.70		
bis(2-Chloroethoxy)methane	.54411	.54835	.78		
2,4-Dichlorophenol	.34085	.34399	.92	*	
1,2,4-Trichlorobenzene	.34396	.42274	22.90		
Naphthalene	1.05928	1.13092	6.76		
4-Chloroaniline		.55768	6.45		
Hexachiorobutadiene		.25357	26.99	*	
4-Chloro-3-methylphenol	. 42543	. 46058	8.26	*	
2-Methylnaphthalene	.65846	.72369	9.91		
Hexachlorocyclopentadiene	.34469	.46478	34.84		**
2,4,6-Trichlorophenol	.43312		.50	*	
2,4,5-Trichlorophenol	.47589	.56839	19.44		
2-Chloronaphthalene	1.15286	1.19020	3.24		
2-Fluorobiphenyl	1.17244	1.22161	4.19		

RF - Response Factor from daily standard file at 50.00 ng/ul

RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/30/95 Time: 10:15

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	X Diff	ccc	SPCC	
2-Nitroaniline	.56115	.51861	7.58			
Dimethylphthalate		1.62114	1.42			
Acenaphthylene	-	1.98413	2.44			
3-Nitroaniline	.74685	.73415	1.70			
Acenaphthene		1.17291	1.70	*		
2,4-Dinitrophenol	.30476	. 29320	3.79		**	
4-Nitrophenol	.47728	.41252	13.57		**	
Dibenzofuran	1.84025	1.91479	4.05			
2,6-Dinitrotoluene	.41593	.40641	2.29			
2,4-Dinitrotoluene	.63725	.63428	.47			
Diethylphthalate	1.48656	1.52092	2.31			
4-Chlorophenyl-phenylether	.76081	.74583	1.97			
Fluorene	1.22830	1.29015	5.04			
4-Nitroaniline	.59631	.56504	5.24			
4,6-Dinitro-2-methylphenol	.20574	.22814	10.89			
N-Nitrosodiphenylamine	.50773	.52100	2.62	*		
2,4,6-Tribromophenol	.17241	.17573	1.92			
4-Bromophenyl-phenylether	.26103	.26972	3.33			
Hexachlorobenzene	.32717	.37776	15.46			
Pentachlorophenol	. 23825	. 26044	9.32	*		
Phenanthrene	1.10818	1.22965	10.96			
Anthracene	1.16825	1.16126	.60			
Carbazole		1.12322	4.38			
Di-n-butylphthalate	1.69660	1.73809	2.45			
Fluoranthene	1.39808	1.55282	11.07	*		
Pyrene		1.50998	2.08			
Terphenyl-d14		1.22594	5.34			
Butylbenzylphthalate		. 93655	3.53			
3,3'-Dichlorobenzidine			5.61			
Benzo(a)anthracene		1.58749	.83			
Bis(2-Ethylhexyl)phthalate		1.13984	4.76			
Chrysene		1.32579	.06			
Di-n-octylphthalate		1.91119	6.98	*		
Benzo(b)fluoranthene		1.46600	1.99			
Benzo(k)fluoranthene		1.28834	19.97			
Benzo(a)pyrene		1.34918	4.33	*		
Indeno(1,2,3-cd)pyrene		1.52596	10.18			
Dibenzo(a,h)anthracene		1.06821	14.21			
Benzo(g,h,i)perylene	1.23341	1.38885	12.60			
DF		 	1			

- Response Factor from daily standard file at 50.00 ng/ul

RF RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/31/95

Time: 09:10

Laboratory ID: >PD021 Instrument ID: 888632

Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	% Diff	ССС	SPCC
2-Fluorophenol	1 14925	1.14648	.24		
Phenol-d5		1.97690	11.89		
Phenol		1.81058	15.17	*	
bis(-2-Chloroethyl)Ether		1.77330	9.77		
2-Chlorophenol		1.59342	6.79		
1,3-Dichlorobenzene		1.59605	5.51		
1,4-Dichlorobenzene	1.55768	1.70840	9.68	*	
Benzyl alcohol	1.58553	1.61039	1.57		
1,2-Dichlorobenzene	1.39231	1.56561	12.45		
2-Methylphenol	1.34514	1.48279	10.23		
bis(2-Chloroisopropyl)ether	3.77777	3.96332	4.91		
4-Methylphenol	1.32191	1.53198	15.89		
N-Nitroso-Di-n-propylamine	1.99373	1.97419	.98		**
Hexachioroethane	.49991	.62472	24.97		
Nitrobenzene-d5	. 43070	.46846	8.77		
Nitrobenzene	.47397	.52176	10.08		
Isophorone	.98795	.94999	3.84		
2-Nitrophenol	. 24923		2.48	*	
2,4-Dimethylphenol	.33290		2.45		
bis(2-Chloroethoxy)methane		.58099	6.78		
2,4-Dichlorophenol	.34085		11.21	*	
1,2,4-Trichlorobenzene	.34396		18.32		
Naphthalene		1.14423	8.02		
4-Chloroaniline	.52389		10.82		
Hexachlorobutadiene	.19968		34.48	*	
4-Chloro-3-methylphenol	.42543		6.44	*	
2-Methylnaphthalene	.65846		19.90		
Hexachlorocyclopentadiene	.34469		20.72	_	**
2,4,6-Trichtorophenol	.43312		.52	*	
2,4,5-Trichlorophenol	.47589		18.34		
2-Chloronaphthalene		1.19866	3.97		
2-Fluorobiphenyl	1.1/244	1.26700	8.06		

RF - Response Factor from daily standard file at 50.00 ng/ul

RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 01/31/95

Time: 09:10

Laboratory ID: >PD021 Instrument ID: 888632

Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	%Diff	CCC	SPCC
2-Nitrosniline	.56115	.57639	2.72		
Dimethylphthalate	1.59849	1.61186	.84		
Acenaphthylene	1.93695	2.03894	5.27		
3-Nitroaniline	.74685	. 78791	5.50		
Acenaphthene	1.19323	1.24582	4.41	*	
2,4-Dinitrophenol	.30476	.23390	23,25		**
4-Nitrophenol	.47728	.41970	12.06		**
Dibenzofuran	1.84025		9.41		
2,6-Dinitrotoluene	.41593	.35221	15,32		
2,4-Dinitrotoluene	.63725	.56946	10.64		
Diethylphthalate	1.48656	1.43015	3.79		
4-Chlorophenyl-phenylether	.76081	.75395	.90		
Fluorene	1.22830		5.36		
4-Nitroaniline	.59631		1.51		
4,6-Dinitro-2-methylphenol	.20574				
N-Nitrosodiphenylamine	.50773			*	
2,4,6-Tribromophenol		. 16515	4.21		
4-Bromophenyl-phenylether		.26797	2.66		
Hexach Lorobenzene	.32717		10.33		
Pentachlorophenol	.23825		3.71	*	
Phenanthrene		1.20302			
Anthracene	1.16825		9.12		
Carbazole	1.17462		7.46		
Di-n-butylphthalate		1.73879	2.49		
Fluoranthene	1.39808		7.48	*	
Pyrene	1,54199		2.28		
Terphenyl-d14	1.16385		.22		
Butylbenzylphthalate	.90464		13.60		
3,3'-Dichlorobenzidine	,56015		13.24		
Benzo(a)anthracene	1.57436		5.24		
Bis(2-Ethylhexyl)phthalate	1.19680		1.97		
Chrysene		1.29057	2.60		
Di-n-octylphthalate		2.09868	17.47	*	
Benzo(b)fluoranthene		1.80656	20.78		
Benzo(k)fluoranthene		1.23303			
Benzo(a)pyrene		1.45138		*	
Indeno(1,2,3-cd)pyrene		1.55444			
Dibenzo(a,h)anthracene		1.08532	16.04		
Benzo(g,h,i)perylene		1.41289	14.55		
senzo(g,n, i)per y cene	1.23341	1.41209	14.33		

RF - Response Factor from daily standard file at 50.00 ng/ul

RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 02/01/95 Time: 10:31

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	XD iff	ccc	SPCC
2-Fiuorophenol	1 14925	1.08586	5.52		
Phenol-d5		1.76498	.10		
Phenol		2.03069	4.85	*	
bis(-2-Chloroethyl)Ether		1.74998	8.33		
2-Chlorophenol		1.57731	5.71		
1,3-Dichlorobenzene		1.60004	5.78		
1,4-Dichlorobenzene	1.55768	1.66518	6.90	*	
Benzyl alcohol	1.58553	1.66734	5.16		
1,2-Dichlorobenzene	1.39231	1.44442	3.74		
2-Methylphenol	1.34514	1.45549	8.20		
bis(2-Chloroisopropyl)ether	3.77777	4.04862	7.17		
4-Methylphenol	1.32191	1.43272	8.38		
N-Nitroso-Di-n-propylamine	1.99373	2.12021	6.34		**
Hexachloroethane		.57768	15.56		
Nitrobenzene-d5	.43070	.44100	2.39		
Nitrobenzene	.47397		4.51		
Isophorone	.987 9 5	.95994	2.84		
2-Nitrophenol	. 24923	. 24054	3.49	*	
2,4-Dimethylphenol	.33290	.33119	.51		
bis(2-Chloroethoxy)methane		.57521	5.72		
2,4-Dichlorophenol		.37785		*	
1,2,4-Trichlorobenzene	.34396		11.50		
Naphthalene	1.05928		6.11		
4-Chloroaniline		.55531	6.00		
Hexach Lorobut adiene	. 19968		38. <i>7</i> 5	*	
4-Chloro-3-methylphenol	.42543		3.67	*	
2-Methylnaphthalene	.65846		14.52		
Hexachlorocyclopentadiene	.34469		11.22		**
2,4,6-Trichlorophenol	.43312		4.24	*	
2,4,5-Trichlorophenol	.47589				
2-Chloronaphthalene	1.15286		6.02		
2-Fluorobiphenyl	1.17244	1.26025	7.49		

RF - Response Factor from daily standard file at 50.00 ng/ul

RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**) RF

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 02/01/95

Time: 10:31

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	X Diff	ccc	SPCC
2-Nitroaniline	.56115	.58069	3.48		
Dimethylphthalate	1.59849	1.68718	5.55		
Acenaphthylene	1.93695	2.09265	8.04		
3-Nitroaniline	. 74685	.82760	10.81		
Acenaphthene	1.19323	1.17760	1.31	*	
2,4-Dinitrophenol	.30476	.20610	32.37		**
4-Nitrophenol	.47728	.44051	7.71		**
Dibenzofuran	1.84025	1.82085	1.05		
2,6-Dinitrotoluene	.41593	.40379	2.92		
2,4-Dinitrotoluene	.63725	.61157	4.03		
Diethylphthalate	1.48656	1.51714	2.06		
4-Chlorophenyl-phenylether	.76081	.80843	6.26		
fluorene	1.22830	1.18262	3.72		
4-Nitroaniline	.59631	.62537	4.87		
4,6-Dinitro-2-methylphenol	.20574	.18640	9.40		
N-Nitrosodiphenylamine	.50773	.54509	7.36	*	
2,4,6-Tribromophenol	.17241	. 17665	2.46		
4-Bromophenyl-phenylether	.26103		9.56		
Hexach Lorobenzene	.32717	.34301	4.84		
Pentachlorophenol	.23825	.22768	4.44	*	
Phenanthrene	1.10818	1.26553	14.20		
Anthracene	1.16825	1.28609	10.09		
Carbazole	1.17462	1.24436	5.94		
Di-n-butylphthalate		1.79358	5.72		
Fluoranthene	1.39808	1.57019	12.31	*	
Pyrene	1.54199	1.38891	9.93		
Terphenyl-d14	1.16385	1.05838	9.06		
Butylbenzylphthalate	.90464	.77666	14.15		
3,3'-Dichlorobenzidine	.56015	.52666	5.98		
Benzo(a)anthracene	1.57436	1.36025	13.60		
Bis(2-Ethylhexyl)phthalate	1.19680	1.10001	8.09		
Chrysene	1.32497	1.20796	8.83		
Di-n-octylphthalate	1.78652	2.00631	12.30	*	
Benzo(b)fluoranthene	1.49577	1.53461	2.60		
Benzo(k)fluoranthene	1.07387	1.31787	22.72		
Benzo(a)pyrene	1.29324	-	14.47	*	
Indeno(1,2,3-cd)pyrene	1.38492		8.49		
Dibenzo(a,h)anthracene		1.00249	7.19		
Benzo(g,h,i)perylene	1.23341		9.06		
	·				

- Response factor from daily standard file at 50.00 ng/ul

RF - Average Response Factor from Initial Calibration form VI %Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 02/08/95

Time: 10:29
Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	XD iff	CCC	SPCC
2-Fluorophenol	1.14925	1.17310	2.08		
Phenol-d5		1.77192	.29		
Phenol		1.86812	12.47	*	
bis(-2-Chloroethyl)Ether	1.61547	1.83714	13.72		
2-Chlorophenol	1.49216	1.57134	5.31		
1,3-Dichlorobenzene	1.51267	1.62652	7.53		
1,4-Dichlorobenzene	1.55768	1.60284	2.90	*	
Benzyl alcohol	1.58553	1.56492	1.30		
1,2-Dichlorobenzene	1.39231	1.39773	.39		
2-Methylphenol	1.34514	1.36788			
bis(2-Chloroisopropyl)ether	3.77777	3.70262	1.99		
4-Methylphenol		1.43650	8.67		
N-Nitroso-Di-n-propylamine	1.99373	1.87177	6.12		**
Hexachloroethane	.49991		18.52		
Nitrobenzene-d5	.43070		7.04		
Nitrobenzene	.47397		9.71		
I sophorone	.98795		.55		
2-Nitrophenol	. 24923	.21218	14.87	*	
2,4-Dimethylphenol	.33290		3.27		
bis(2-Chloroethoxy)methane		.58024	6.64		
2,4-Dichlorophenol	.34085		10.23	*	
1,2,4-Trichlorobenzene	.34396	.40331	17.25		
Naph tha lene	1.05928		8.79		
4-Chloroaniline	.52389		6.86		
Hexach Lorobutadiene	. 19968		22.70	*	
4-Chloro-3-methylphenol	.42543		9.97	*	
2-Methylnaphthalene	.65846		7.33		
Hexachlorocyclopentadiene		.43171	25.25		**
2,4,6-Trichlorophenol		.44355		*	
2,4,5-Trichlorophenol		.53020			
2-Chloronaphthalene		1.23594	7.21		
2-Fluorobiphenyl	1.1/244	1.29828	10.73		

- Response Factor from daily standard file at 50.00 ng/ul

RF - Average Response Factor from Initial Calibration Form V1 %Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 02/08/95

Time: 10:29

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	X Diff	ccc	SPCC
2-Nitroaniline	.56115	.59087	5.30		
Dimethylphthalate	1.59849	1.71422	7.24		
Acenaphthylene	1.93695	2.15952	11.49		
3-Nitroaniline	.74685	.85054	13.88		
Acenaphthene	1.19323	1.24739	4.54	*	
2,4-Dinitrophenol	.30476	.27378	10.17		**
4-Nitrophenol	.47728	.46081	3.45		**
Dibenzofuran	1.84025	1.98320	7.77		
2,6-Dinitrotoluene	.41593	.45855	10.25		
2,4-Dinitrotoluene	.63725	.63862	.21		
Diethylphthalate	1.48656	1.42611	4,07		
4-Chlorophenyl-phenylether	.76081	.77920	2,42		
Fluorene	1.22830	1.32750	8.08		
4-Nitroaniline	.59631	.60793	1.95		
4,6-Dinitro-2-methylphenol	.20574	.21140	2.75		
N-Nitrosodiphenylamine	.50773	.58384	14.99	*	
2,4,6-Tribromophenol		.16740	2.91		
4-Bromophenyl-phenylether	.26103	. 26985	3.38		
Hexach Lorobenzene	.32717	.34086	4.18		
Pentachlorophenol	.23825		7.49	*	
Phenanthrene	1.10818		4.67		
Anthracene	1.16825		.42		
Carbazole	1.17462		.95		
Di-n-butylphthalate	1.69660		8.12		
Fluoranthene	1.39808		5.74	*	
Pyrene	1.54199		3.36		
Terphenyl-d14	1.16385		6.60		
Butylbenzylphthalate	.90464		3.04		
3,31-Dichlorobenzidine	.56015		6,23		
Benzo(a)anthracene	1,57436		2.25		
Bis(2-Ethylhexyl)phthalate	1.19680		5.98		
Chrysene	1.32497		.77		
Di-n-octylphthalate	1,78652		8.20	*	
Benzo(b)fluoranthene	1,49577		6.24		
Benzo(k)fluoranthene	1,07387		11.83		
Benzo(a)pyrene	1.29324		2.34	*	
Indeno(1,2,3-cd)pyrene	1.38492		4.06		
Dibenzo(a,h)anthracene	.93528		4.89		
Benzo(g,h,i)perylene	1.23341		8.73		
	1,23341				
05 0 5 6					

RF - Response Factor from daily standard file at 50.00 ng/ul

- Average Response Factor from Initial Calibration Form VI

%Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)
SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 02/24/95

Time: 10:38

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	XD iff	CCC	SPCC
2-Fluorophenol	1 14025	1.11077	3.35		
Phenol-d5		1.78025	.76		
Phenol		1.99234		*	
bis(-2-Chloroethyl)Ether		1.78228	10.33		
2-Chlorophenol	1.49216	1.49923	.47		
1,3-Dichlorobenzene	1.51267	1.74437	15.32		
1,4-Dichlorobenzene	1.55768	1.70036	9.16	*	
Benzyl alcohol	1.58553	1.51099	4.70		
1,2-Dichlorobenzene	1.39231	1.75825	26.28		
2-Methylphenol	1.34514	1.52546	13.41		
bis(2-Chloroisopropyl)ether	3.77777	4.27058	13.04		
4-Methylphenol	1.32191	1.73935	31.58		
N-Nitroso-Di-n-propylamine	1.99373	2.17541	9.11		**
Hexachloroethane	.4999 1		31.48		
Nitrobenzene-d5	.43070	. 45436	5.49		
Nitrobenzene	.47397	.47849	.95		
Isophorone	. 98795	.97719	1.09		
2-Nitrophenol	. 24923		3.24	*	
2,4-Dimethylphenol		.30837	7.37		
bis(2-Chloroethoxy)methane		.54747	.62		
2,4-Dichlorophenol		.31339	8.06	*	
1,2,4-Trichlorobenzene	.34396		20.24		
Naphthalene		1.19560	12.87		
4-Chloroaniline	.52389		3.19		
Hexachlorobutadiene	. 19968		21.16	*	
4-Chloro-3-methylphenol	.42543		1.46	*	
2-Methylnaphthalene	.65846		16.03		
Hexachlorocyclopentadiene	.34469		18.84		**
2,4,6-Trichlorophenol		.40334	6.87	*	
2,4,5-Trichlorophenol	.47589		. 18		
2-Chloronaphthalene		1.16981			
2-Fluorobiphenyl	1.17244	1.15068	1.86		

- Response Factor from daily standard file at 50.00 ng/ul RF

RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve CCC - Calibration Check Compounds (*) SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 02/24/95

Time: 10:38

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	X Diff	CCC	SPCC
2-Nitroaniline	.56115	.51753	7.77		
Dimethylphthalate		1.67345	4.69		
Acenaphthylene		2.18496			
3-Nitroaniline	.74685		4.46		
Acenaphthene		1.22196	2.41	*	
2,4-Dinitrophenol		.23414	23.17		**
4-Nitrophenol	.47728		13.69		**
Dibenzofuran		1.78510	3.00		
2,6-Dinitrotoluene		.42342			
2,4-Dinitrotoluene	.63725	.61659	3.24		
Diethylphthalate	1.48656	1.54854	4.17		
4-Chlorophenyl-phenylether	.76081	.79115	3.99		
Fluorene	1.22830	1.28462	4.59		
4-Nitroaniline	.59631	.42477	28.77		
4,6-Dinitro-2-methylphenol	. 20574	.21185	2.97		
N-Nitrosodiphenylamine	.50773	.56765	11.80	*	
2,4,6-Tribromophenol	.17241				
4-Bromophenyl-phenylether	.26103	.26116	.05		
Hexach Lorobenzene	.32717	.34767	6.27		
Pentachlorophenol		.24062	.99	*	
Phenanthrene		1.28922	16.34		
Anthracene	1.16825	1.10498	5.42		
Carbazole		1.20880	2.91		
Di-n-butylphthalate		1.87825	10.71		
fluoranthene	1.39808	1.58180	13.14	*	
Pyrene		1.55345	.74		
Terphenyl-d14		1.03407	11.15		
Butylbenzylphthalate	.90464		4.17		
3,3'-Dichlorobenzidine	.56015		12.18		
Benzo(a)anthracene		1.38007	12.34		
Bis(2-Ethylhexyl)phthalate		1.27244	6.32		
Chrysene		1.28708	2.86		
Di-n-octylphthalate		2.34797	31.43	*	
Benzo(b)fluoranthene		1.84985	23.67		
Benzo(k)fluoranthene		1.06092	1.21		
Benzo(a)pyrene		1.41909		*	
Indeno(1,2,3-cd)pyrene		1.66056			
Dibenzo(a,h)anthracene		1.04955			
Benzo(g,h,i)perylene	1.23341	1.47549	19.63		

RF - Response Factor from daily standard file at 50.00 ng/ul

- Average Response Factor from Initial Calibration Form VI

2 2 34.5

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 02/27/95

Time: 15:44

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	2Diff	ccc	SPCC
2-Fluorophenol	1.14925	1.06731	7,13		
Phenol-d5	1.76675	1.68367	4.70		
Phenol	2.13426	1.86727	12.51	*	
bis(-2-Chloroethyl)Ether	1.61547	1.73528	7.42		
2-Chlorophenol	1.49216	1.42290	4.64		
1,3-Dichlorobenzene	1.51267	1.51627	.24		
1,4-Dichlorobenzene	1.55768	1.62937	4.60	*	
Benzyl alcohol	1.58553	1.70536	7.56		
1,2-Dichlorobenzene	1.39231	1.57515	13.13		
2-Methylphenol	1.34514	1.26173	6.20		
bis(2-Chloroisopropyl)ether		3.67622			
4-Methylphenol	1.32191				
N-Nitroso-Di-n-propylamine		2.08075			**
Hexachloroethane		.62742			
Nitrobenzene-d5		.47644	10.62		
Nitrobenzene	.47397		1.53		
Isophorone		1.02135	3.38		
2-Nitrophenol	. 24923		1.22	*	
2,4-Dimethylphenol	.33290		1.62		
bis(2-Chloroethoxy)methane	.54411		2.12		
2,4-Dichtorophenot	.34085		2.14	*	
1,2,4-Trichlorobenzene	.34396		9.20		
Naphthalene		1.15502	9.04		
4-Chloroaniline	.52389	_	-		
Hexachlorobutadiene	. 19968			*	
4-Chloro-3-methylphenol	.42543		1.64	*	
2-Methylnaphthalene		.73822			
Hexachlorocyclopentadiene		.42271	22.63		**
2,4,6-Trichlorophenol	.43312		1.98	*	
2,4,5-Trichlorophenol	.47589		4.32		
2-Chloronaphthalene		1.28790			
2-Fluorobiphenyl	1.17244	1.24757	6.41		

RF - Response Factor from daily standard file at 50.00 ng/ul

RF - Average Response Factor from Initial Calibration Form VI %Diff - % Difference from original average or curve

CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)

Table 2.26 (Cont) Results of the Continuing Calibration WA # 0-098 Pier Drum

Calibration Date: 02/27/95

Time: 15:44

Instrument ID: 888632 Initial Calibration Date: 11/04/94

Minimum RF for SPCC is 0.05

Maximum % Diff for CCC is 25%

Compound	RF	RF	X Diff	CCC	SPCC
2-Nitroaniline	.56115	.54620	2.66		
Dimethylphthalate	1.59849	1.68315	5.30		
Acenaphthylene	1.93695	1.99336	2.91		
3-Nitroaniline	.74685	. 78099	4.57		
Acenaphthene	1.19323	1.25573	5.24	*	
2,4-Dinitrophenol	.30476	.28857	5.31		**
4-Nitrophenol	.47728	. 48302	1.20		**
Dibenzofuran	1.84025	1.85246	.66		
2,6-Dinitrotoluene	.41593	.38472	7.50		
2,4-Dinitrotoluene	.63725	.68164	6.97		
Diethylphthalate	1.48656	1.66186	11.79		
4-Chlorophenyl-phenylether	.76081	.84581	11.17		
Fluorene	1.22830	1.58742	29.24		
4-Nitroaniline	.59631	.52456	12.03		
4,6-Dinitro-2-methylphenol	.20574	.21662			
N-Nitrosodiphenylamine	.50773			*	
2,4,6-Tribromophenol		. 16254	5.73		
4-Bromophenyl-phenylether	.26103	.25402	2,68		
Hexachlorobenzene	.32717		.31		
Pentachlorophenol	.23825	.22684	4.79	*	
Phenanthrene	1.10818		18.41		
Anthracene	1.16825		8.97		
Carbazole	1.17462		.37		
Di-n-butylphthalate	1.69660		5.41		
Fluoranthene	1.39808		3.65	*	
Pyrene	1,54199		11.07		
Terphenyl-d14	1.16385		10.98		
Butylbenzylphthalate	.90464		10.46		
3,3'-Dichlorobenzidine		.49881	10.95		
Benzo(a)anthracene	1.57436		15.58		
Bis(2-Ethylhexyl)phthalate	1.19680		9.16		
Chrysene	1.32497		15.22		
Di-n-octylphthalate	1.78652		22.86	*	
Benzo(b)fluoranthene	1.49577		4.00		
Benzo(k)fluoranthene	1.07387		33.55		
Benzo(a)pyrene	1.29324			*	
Indeno(1,2,3-cd)pyrene	1.38492				
Dibenzo(a,h)anthracene		1.07056			
Benzo(g,h,i)perylene	1.23341		22.15		

- Response Factor from daily standard file at 50.00 ng/ul RF

RF - Average Response Factor from Initial Calibration Form VI

**Diff - * Difference from original average or curve CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)

Table 2.27 Correlation of Samples and Calibrations

Samples	Initial Calibration	Continuing Calibration
WBLK10199402, PE10199402, SBLK10189402, PE10189402	11/4/94	_
A-C 00759 MS, A-C 00759 MSD, WBLK012695, A 00763, A-C 00764, A-C 00764 MS, A-C 00764 MSD, WBLK011995, B 00525, B 00526, BE 00527,	11/4/94	1/27/95
BE 00527 MS, BE 00527 MSD, WBLK011895 CD00636, CD 00638, CD 00641, CD 00635, CD 00639, CD 00640, CD 00640 MS, CD 00640 MSD, CD 00634, CD 00637	11/4/94	1/30/95
Soil Blank 011395, A 00642, A 00643, A 00644 A 00544 MS, A 00644 MSD	11/4/94	1/31/95
CD 00637	11/4/94	2/1/95
00913 Bott	11/4/94	2/8/95
WBLK 021495, WBLK 021495 MS, WBLK 021495 MSD 00912, 00913, 00914	11/4/94	2/24/95
00912 (10x), 00913 (10x), 00914 (10x)	11/4/94	2/27/95

Table 2.28 Results of the TCLP Surrogate Recoveries for the Drum Solid Samples WA # 0-098 Pier Drum

Sample ID	S1 (NBZ)	S2 (FBP)	S3 (TPH)	S4 (PHL)	S5 (2FP)	S6 (DCB)	Total Out
SVBL	96	80	83	92	90	29	0
Blank MS	83	81	87	83	78	33	0
00915 MS	81	77	90	28	NC*	30	0
00915 MSD	97	91	94	22	NC *	87	0
00915 MSD (DL)	62	83	107	24	40	52	0
00911	111	60	106	38	11 *	52	1
00915	105	88	109	35	41	51	0
00916	105	92	104	41	NC *	55	0
00916 (DL)	62	94	100	22	38	53	0

* denotes a value that exceeds acceptable QC limits or was not calculated due to interferences

OC Limits

QC LIMITS
35-114
43-116
33-141
10-110
21-110
10-123

Table 2.29 Results of the TCLP Internal Standard Areas for the Drum Solid Samples WA # 0-098 Pier Drum

ISTD 2 ISTD 3 ISTD 4 ISTD 5 ISTD 6 ISTD 1 Total Sample ID **AREA AREA AREA** AREA AREA AREA Out 50 ppb Std SVBL Blank MS 00915 MS 147786 * 50 ppb Std 00916 (DL) O 50 ppb Std 00915 MSD 50 ppb Std 00915 MSD (DL)

The acceptable QC limts are 50 - 200% of the appropriate standard

ISTD 1 denotes 1,4-dichlorobenzene-d4

ISTD 2 denotes naphthalene-d8

ISTD 3 denotes acenaphthene-d10

ISTD 4 denotes phenanthrene-d10

ISTD 5 denotes chrysene-d12

ISTD 6 denotes perylene-d12

Table 2.30 Results of the Blank Spike Analysis for the TCLP BNA Analysis WA # 0-098 Pier Drum

	Spike	Sample	BS		QC Limits
Parameter	Added	Conc.	Recov.	%	
	(u g/L)	(ug/L)	(ug/L)	Recovery	Rec.
1,2,4-Trichlorobenzene	100	ND	48	48	39-98
Acenapthene	100	ND	79	79	46-118
2,4-Dinitrotoluene	100	ND	42	42	24-96
Pyrene	100	ND	104	104	26-127
N-Nitroso-di-n-propylamine	100	ND	61	61	41-116
1,4-Dichlorobenzene	100	ND	56	56	36-97
Pentachiorophenol	100	ND	73	73	9-103
Phenol	100	ND	143	143	12-110
2-Chlorophenol	100	ND	150	150	27-123
4-Chloro-3-methylphenol	100	ND	127	127	23-97
4-Nitrophenol	100	ND	67	67	10-80

Table 2.31 Results of the MS/MSD Analysis for the TCLP BNA Analysis of the Drum Solid Samples WA # 0-098 Pier Drum

Sample ID: 00915

Danamatan	Spike Added	Sample Conc.	MS Recov.	%	MSD Recov.	<u>-</u> -			QC mits
Parameter	(ug/L)	(ug/L)	(ug/L)	Recovery	(ug/L)	Recovery	RPD	RPD	Rec.
2-Chlorophenol	444	ND	302	68	264	59	14	40	27-123
Phenol	444	17	109	21	120	23	10	42	12-110
1,4 - Dichlorobenzene	222	ND	87	39	128	58	39 *	28	36-97
N-Nitroso-di-n-propylamine	222	ND	140	63	146	66	4	38	41-116
1,2,4-Trichlorobenzene	222	ND	79	36 *	144	65	58 *	28	39-98
4-Chloro-3-methylphenol	444	ND	289	6 5	294	66	2	42	23-97
Acenapthene	222	ND	162	73	154	69	5	31	46-118
4 – Nitrophenol	444	ND	54	12	92	21	52 *	50	10-80
2,4-Dinitrotoluene	222	ND	93	42	170	77	59 *	38	24-96
Pentachlorophenol	444	ND	235	53	369	83	44	50	9-103
Pyrene	222	ND	211	95	105	47	66 *	31	_26-127

7 2 1433

QA/QC FOR PESTICIDES/PCBs

Each sample was spiked with a solution of tetrachloro-m-xylene and decachlorobiphenyl as surrogates; percent recoveries for the water samples, listed in Table 2.32, ranged from 20 to 83. Nine out of twenty-two values were within the advisory QC limits.

Sample 00640 was chosen for the matrix spike/matrix spike duplicate (MS/MSD) analyses for the water samples. The percent recoveries, ranging from 50 to 88, are listed in Table 2.33. All twelve percent recoveries are within acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.33, ranged from 0 (zero) to 9 and all six values were within the acceptable QC limits.

The surrogate percent recoveries for the soil samples, listed in Table 2.34, ranged from 26 to 132. Sixteen out of eighteen values were within the advisory QC limits.

Samples 00642 and 00643 were chosen for the MS/MSD analyses for the soil samples. The percent recoveries, ranging from 22 to 152, are listed in Table 2.35. Twenty out of twenty-four percent recoveries were within the acceptable QC limits. The RPDs, also listed in Table 2.35, ranged from 2 to 131, and five out of twelve values were within the acceptable QC limits.

The surrogate percent recoveries for the TCLP samples, listed in Table 2.36, ranged from 19 to 95. Five out of twelve values were within the advisory QC limits.

Sample CF 00527 was chosen for the MS/MSD analyses for the TCLP samples. The percent recoveries, ranging from 64 to 109, are listed in Table 2.37. All twelve percent recoveries were within the acceptable QC limits. The RPDs, also listed in Table 2.37, ranged from 0 (zero) to 11, and all six values were within the acceptable QC limits.

The surrogate percent recoveries for the drum liquid samples, listed in Table 2.38, ranged from 14 to 76. One out of nine values were within the advisory QC limits. Five other values were not calculated due to matrix interference.

Sample WBLK020795 was chosen for the MS/MSD analyses for the drum liquid samples. The percent recoveries, ranging from 56 to 88, are listed in Table 2.39. All twelve percent recoveries were within the acceptable QC limits. The RPDs, also listed in Table 2.39, ranged from 0 (zero) to 13, and all six values were within the acceptable QC limits.

The surrogate percent recoveries for the PCB analysis of the TCLP extracts of the drum solids sample, listed in Table 2.40A, ranged from 26 to 87. All eight values were within the advisory QC limits. Two values were from diluted samples and the surrogate was not recovered.

The surrogate percent recoveries for the pesticide analysis of the TCLP extracts of the drum solids sample, listed in Table 2.40B, ranged from 42 to 121. All ten values were within the advisory QC limits.

2 2 3433

The results of the blank spike, listed in Table 2.41, ranged from 52 to 100. Six out of seven percent recoveries were within the acceptable QC limits.

Sample 00911 was chosen for the MS/MSD for the pesticide analyses for the TCLP extracts of the drum solid samples. The percent recoveries, ranging from 14 to 182, are listed in Table 2.42. Six out of twelve percent recoveries were within the acceptable QC limits. The RPDs, also listed in Table 2.42, ranged from 44 to 124, and none of the six values were within the acceptable QC limits.

Sample 00916 was chosen for the MS/MSD for the PCB analyses for the TCLP extracts of the drum solid samples. The percent recoveries were 88 and 125 and are also listed in Table 2.42. Both percent recoveries were within the acceptable QC limits. The RPD, also listed in Table 2.42, was 35, and it was within the acceptable QC limits.

Table 2.32 Results of the Surrogate Recoveries in Water WA # 0-098 Pier Drum

	Percent F	Recovery
Sample ID	TCMX	DCBP
WBLK011795	29 *	62
EF 00636	51 *	83
EF 00638	36 *	80
EF 00641	33 *	64
EF 00635	41 *	69
EF 00639	20 *	58 *
EF 00640	25 *	64
EF 00640MS	43 *	70
EF 00640MSD	38 *	63
EF 00634	28 *	46 *
EF 00637	33 *	70

TCMX denotesTetrachloro – m – xylene DCBP denotes Decachlorobiphenyl

	Advisory
	QC
	Limits
TCMX	60-150
DCBP	60-150

Table 2.33 Results of the MS/MSD Analysis for Pest/PCB in Water WA # 0-098 Pier Drum

Sample ID: EF 00640

Compound	Sample Conc	MS Spike Added	M S Conc	MS %	MSD Spike Added	M S D Conc	MSD %)	QC Lir %	nit
·	(µg/L)	(ug/L)	(ug/L)	Rec	(µg/L)	(n8/r)	Rec	RPD	Rec	RPD
g-BHC	ND	0.20	0.14	71	0.20	0.15	76	7	46-127	50
Heptachlor	ND	0.20	0.11	55	0.20	0.12	61	9	35-130	31
Aldrin	ND	0.20	0.10	50	0.20	0.10	50	0	34-132	43
Dieldrin	ND	0.40	0.27	68	0.40	0.29	73	7	31-134	38
Endrin	ND	0.40	0.30	76	0.40	0.32	81	6	42-139	45
p,p'-DDT	ND	0.40	0.34	86	0.40	0.35	88	3	23-134	50

Table 2.34 Results of the Surrogate Recoveries in Soil WA # 0-098 Pier Drum

Sample iD	Percent	Recovery
SBLK011395 A 00643 A 00643MS A 00643MSD A 00644 SBLK020795 A 00642	TCMX	DCBP
SBLK011395	117	110
A 00643	82	31 *
A 00643MS	94	126
A 00643MSD	68	26 *
A 00644	69	78
SBLK020795	98	132
A 00642	105	89
A 00642 MS	84	88
A 00642 MSD	98	101

TCMX denotesTetrachloro-m-xylene DCBP denotes Decachlorobiphenyl

	Advisor
	QC
	Limits
TCMX	60-150
DCBP	60~150

098\DEL\A > 2E.

00170

Table 2.35 Results of the MS/MSD Analysis for Pest/PCB in Soil WA # 0-098 Pier Drum

Sample ID: A 00643

Compound	Sample Conc		M S Conc	M S _%	MSD Spike Added	M S D Conc	MSD _%		QC Lir %	
	(µg/Kg)	(µg/Kg)	(µg/Kg)	Rec	(µg/Kg)	(µg/Kg)	Rec	RPD	Rec	RPD
g-BHC	ND	25	28	112	25	18	72	43	46-127	50
Heptachlor	ND	25	26	104	25	15	60	54 *	35-130	31
Aldrin	ND	25	26	104	25	5	22 1	131 *	34-132	43
Dieldrin	ND	50	60	120	50	33	66	58 *	31-134	38
Endrin	ND	50	70	140 *	50	37	74	62 *	42-139	45
p.p'-DDT	ND	50	76	152 *	50	33	66	79 *	23-134	50

Sample ID: A 00642

Compound	Sample Conc	MS Spike Added	M S Conc	MS %	MSD Spike Added	M S D Conc	MSD %	1	QC Lir %	nit
·	(μg/Kg)	(µg/Kg)	(µg/Kg)	Rec	(µg/Kg)	(µg/Kg)	Rec	RPD	Rec	RPD
g-BHC	ND	25.6	18	70	25.6	21	82	15	46-127	50
Heptachlor	ND	25.6	20	78	25.6	22	86	10	35-130	31
Aldrin	ND	25.6	37	145 *	25.6	25	98	39	34-132	43
Dieldrin	ND	51.3	41	80	51.3	44	86	7	31-134	38
Endrin	ND	51.3	42	82	51.3	43	84	2	42-139	45
p,p'-DDT	ND	51.3	36	70	51.3	44	86	20	23-134	50

Table 2.36 Results of the Surrogate Recoveries in TCLP WA # 0-098 Pier Drum

	Z		
Sample ID	TCMX	DCBP	
WBLK011995	19 *	49 *	
C 00525	38 *	85	
C 00526	25 *	78	
C 00527	49 *	88	
C 00527MS	47 *	86	
C 00527MSD	48 *	95	

TCMX denotesTetrachioro –m –xylene DCBP denotes Decachiorobiphenyl

	Advisory
	QC
	Limits
TCMX	60-150
DCBP	60-150

Table 2.37 Results of the MS/MSD Analysis for Pest/PCB in TCLP WA # 0-098 Pier Drum

2 11 5439

Sample ID: CF 00527

Compound	Sample Conc (µg/L)	MS Spike Added (µg/L)	M S Conc (µg/L)	MS % Rec	MSD Spike Added (µg/L)	MSD Conc (µg/L)	MSD % Rec	RPD	QC Lin % Rec	nit RPD
g-BHC	ND	0.22	0.18	82	0.22	0.20	91	11	46-127	50
Heptachlor	ND	0.22	0.14	64	0.22	0.15	68	7	35-130	31
Aldrin	ND	0.22	0.15	68	0.22	0.15	68	0	34-132	43
Dieldrin	ND	0.44	0.38	86	0.44	0.39	89	3	31-134	38
Endrin	ND	0.44	0.41	93	0.44	0.43	98	5	42-139	45
p,p'-DDT	ND	0.44	0.44	100	0.44	0.48	109	9	23-134	50

Table 2.38 Results of the Surrogate Recoveries in Drum Liquids WA # 0-098 Pier Drum

WBLK020795	36 *	52 *
WBLK020795 MS	38 *	59 *
WBLK020795 MSD	38 *	76
00912	MI	MI
00913 Top layer	Mi	Mi
00913 Bottom Layer	14 *	MI
00914	23 *	51 *

TCMX denotesTetrachloro –m –xylene DCBP denotes Decachlorobiphenyl

	Advisory
	QC
	Limits
TCMX	60-150
DCBP	60-150

Table 2.39 Results of the MS/MSD Analysis for Pest/PCB in Drum Liquids WA # 0 – 098 Pier Drum

Sample ID: WBLK020795

Compound	Sample Conc	MS Spike Added	M S Conc	MS %	MSD Spike Added	M S D Conc	MSD %)	QC Lir %	nit
	(H g /L)	(ha\r)	(µg/L)	Rec	(h ð /L)	(h8/r)	Rec	RPD	Rec	RPD
g-BHC	ND	0.125	0.090	72	0.125	0.100	80	11	56-123	15
Heptachlor	ND	0.125	0.070	56	0.125	0.080	64	13	40-131	20
Aldrin	ND	0.125	0.070	56	0.125	0.070	56	0	40~120	22
Dieldrin	ND	0.250	0.180	72	0.250	0.200	80	11	52-126	18
Endrin	ND	0.250	0.180	72	0.250	0.200	80	11	56~121	21
p,p'-DDT	ND	0.250	0.200	80	0.250	0.220	88	10	38-127	27

117.

Table 2.40A Results of the Surrogate Recoveries for the Drum Solid Samples WA # 0-098 Pier Drum

2 1 344.2

Sample ID	Percent Recovery of Dibutyl Chlorendate		
Method Blank	87		
Blank Spike	83		
00916 MS	6 5		
00916 MSD	49		
00911	61		
00911 (10x)	D		
00915	46		
00915 (10x)	D		
00916	77		
00916 (10x)	26		

QC Limits

Dibutyi Chiorendate 24-143

Table 2.40B Results of the Surrogate Recoveries for the Drum Solid Samples

WA # 0-098 Pier Drum

Sample ID	Percent Recovery of Dibutyl Chlorendate		
Method Blank	81		
Blank Spike	82		
00911 MS	103		
00911 MSD	42		
00911	121		
00911 (10x)	71		
00915	94		
00915 (10x)	114		
00916	106		
00916 (10x)	76		

QC Limits

Dibutyl Chlorendate 25-154

Table 2.41 Results of the Blank Spike Analysis for the Drum Solid Samples WA # 0-098 Pier Drum

Parameter	Spike Added	Sample Conc. (ug/L)	MS Recov. (ug/L)	%	QC Limits	
	(ug/L)			Recovery	Rec.	
Lindane	0.5	ND	0.39	78	56-123	
Aldrin	0.5	ND	0.37	74	40-120	
Heptachior	0.5	ND	0.34	68	40-131	
Dieldrin	0.5	ND	0.40	80	52-126	
Endrin	0.5	ND	0.26	52 *	56-121	
4,4'-DDT	0.5	ND	0.30	60	38-127	
Arocior1260	2.0	ND	2.0	100	20-150	

Table 2.42 Results of the MS/MSD Analysis for the Drum Solid Samples WA # 0-098 Pier Drum

Sample ID: 00911

	Spike	Sample	MS		MSD				QC
Parameter	Added	Conc.	Recov.	_ %	Recov.	_ %		_	mits
	(ug/L)	(ug/L)	(ug/L)	Recovery	(ug/L)	Recovery	RPD	RPD	Rec.
Lindane	2.50	ND	1.24	49 *	0.55	22 *	76 1	50	56-123
Aldrin	2.50	ND	1.99	79	0.96	40	66 ¹	43	40-120
Heptachlor	2.50	ND	4.14	1 6 6 *	2.65	106	44 1	31	40-131
Dieldrin	2.50	ND	1.50	60	0.35	14 *	124 1	38	52-126
Endrin	2.50	ND	4.55	182 *	2.33	93	65 1	45	56-121
4.4'-DDT	2.50	ND	1.31	52	0.62	25 *	72 1	50	38 - 127
Aroclor 1260**	10	ND	12.5	125	8.8	88	35	45	20-150

^{**} denotes that sample 00916 was used

QA/QC FOR METALS

Quality control standards QCS/ICP-2, ERA-419, QC-19x100, TMMA #1 and TMMA #2 were used to check the accuracy of the calibration curve for the water samples. The percent recoveries ranged from 95 to 109 and all twenty-three percent recoveries were within the 95% confidence limits. The 95% confidence limits for ten values (QC-19x100) are not available. The percent recoveries for the water samples are listed in Table 2.43.

Samples A 00636 and B 00635 were chosen for matrix spike/matrix spike duplicate (MS/MSD) analyses for the water samples. The percent recoveries, listed in Table 2.44, ranged from 65 to 127. Seventy out of seventy-six values were within the acceptable QC limits. The relative percent differences (RPDs), also listed in Table 2.44, ranged from 0 (zero) to 26. Thirty-seven out of thirty-eight values were within the acceptable QC limits.

The results of the blank spike analysis for the water samples are reported in Table 2.45. The percent recoveries ranged from 75 to 119. All twenty-three values were within the acceptable QC limits.

Quality control standards QCS/ICP-2, ERA-419, QC-19x100, TMMA #1 and TMMA #2 were used to check the accuracy of the calibration curve for the soil samples. The percent recoveries ranged from 98 to 127 and twenty-three out of twenty-four percent recoveries were within the 95% confidence limits. The 95% confidence limits for thirteen values (QC-19x100) are not available. The percent recoveries for the soil samples are listed in Table 2.46.

Sample A 00643 was chosen for matrix spike/matrix spike duplicate (MS/MSD) analyses for the soil samples. The percent recoveries, listed in Table 2.47, ranged from 62 to 109. The relative percent differences (RPDs), also listed in Table 2.47, ranged from 0 (zero) to 20. Quality control limits for the percent recoveries and the RPDs are not available.

The results of the blank spike analysis are reported in Table 2.48. The percent recoveries ranged from 82 to 114. Quality control limits for the percent recoveries are not available.

Quality control standards ERA-419, QC-7x100, QC-19x100, and TMMA #1 were used to check the accuracy of the calibration curve for the TCLP samples. The percent recoveries ranged from 95 to 108 and all eight percent recoveries were within the 95% confidence limits. The 95% confidence limits for four values (QC-7x100 and QC-19x100) are not available. The percent recoveries for the TCLP samples are listed in Table 2.49.

Samples A 00527 and A,D 00527 were chosen for MS/MSD analyses for the TCLP samples. The percent recoveries, listed in Table 2.50, ranged from 64 to 105. Fourteen out of sixteen values were within the acceptable QC limits. The RPDs, also listed in Table 2.50, ranged from 0 (zero) to 3. All eight values were within the acceptable QC limits.

The results of the blank spike analysis for the TCLP samples are reported in Table 2.51. The percent recoveries ranged from 97 to 109. All eight values were within the acceptable QC limits.

2 1 2/16

Quality control standards ERA-419, QC-7x100, QC-19x100, TMMA #1 and TMMA #2 were used to check the accuracy of the calibration curve for the drum liquid samples. The percent recoveries ranged from 94 to 121 and all eighteen recoveries were within the 95% confidence limits. The 95% confidence limits for seventeen values (QC-7x100 and QC-19x100) are not available. The percent recoveries for the drum liquid samples are listed in Table 2.52.

The results of the blank spike analysis for the drum liquid samples are reported in Table 2.53. The percent recoveries ranged from 95 to 246. Twenty-one out of twenty-two values were within the acceptable QC limits.

The results of the blank spike analysis for the analysis of the TCLP extracts of the drum solid samples are reported in Table 2.54. The percent recoveries ranged from 81 to 111 and all twenty-three values were within the acceptable QC limits.

Sample 00911 was chosen for MS/MSD analyses for the TCLP extracts of the drum solid samples. The percent recoveries, listed in Table 2.55, ranged from 63 to 118. Twenty-nine out of thirty-four values were within the acceptable QC limits. Ten values were not calculated because the concentration of analyte spiked was less than the concentration of analyte in the sample and two values were not calculated because of matrix interference. The RPDs, also listed in Table 2.55, ranged from 0 (zero) to 10. All seventeen values were within the acceptable QC limits. Five values were not calculated because the concentration of analyte spiked was less than the concentration of analyte in the sample and one value was not calculated because of matrix interference.

2 .. 9477

Table 2.43 Results of the QC Standard Analysis for Water Samples WA # 0-098 Pier Drum

METAL	DATE ANALYZED	QUALITY CONTROL STANDARD	CONC. RECOVERED ug/l	TRUE VALUE ug/l	95 % CONFIDENCE INTERVAL	% RECOVERY
Aluminum		QCS/ICP-2	1059	1000	870 - 1100	106
	01/19/95	ERA 419	220	210	150 – 263	105
Antimony	01/17/95	TMMA #2	109	100	81.65-125.67	109
Arsenic	01/17/95	TMMA #1	49.0	50	41.9-55.9	98
Barium	01/1 9/9 5 01/19/95	QCS/ICP-2 ERA 419	1024 202	1000 204	890 - 1070 167 - 241	102 9 9
Beryllium	01/19/95 01/19/95	QC-19 x100 ERA 419	1016 58	1 00 0 56	N/A 44 - 66	102 104
Cadmium	01/19/95	TMMA #1	5.20	5	4.10-5.83	104
Calcium	01/19/95	QC-19 x100	1005	1000	N/A	, 101
Chromium	01/18/95	TMMA #1	48.4	50	43.91 – 56.15	97
Cobalt	01/19/95 01/19/95	QC-19 x100 ERA 419	1021 110	1 00 0 103	N/A 85 - 122	102 107
Copper	01/19/95 01/19/95	QC-19 x100 ERA 419	1022 79	1000 77	N/A 63 - 91	102 103
Iron	01/19/95 01/19/95	QC-19 x100 ERA 419	1039 131	1000 125	N/A 93 - 147	104 105
Lead	01/19/95	TMMA #1	51.5	50	43.4-56.3	103
Magnesium	101/19/95	QC-19 x100	983	1000	N/A	98
Manganese	01/19/95 01/19/95	QC-19 x100 ERA 419	1016 126	1000 122	N/A 100 – 144	102 103
Mercury	01/20/95	TMMA #1	1.9	2	1.40-2.49	95
Nickel	01/19/95 01/19/95	QC-19 x100 ERA 419	1026 97	1000 92	N/A 75 – 110	103 105
Potassium	01/19/95	QCS/ICP-2	9519	10000	8750 - 11100	95
Selenium	01/17/95	TMMA #1	51	50	39.4-57.4	102
Silver	01/26/95	TMMA #1	10.2	10	8.44-11.8	102
Sodium	01/19/95	QCS/ICP-2	1015	1000	830 – 1470	102
Thallium	01/17/95	TMMA #2	51.4	50	39.9-57.97	103
Vanadium	01/19/95 01/19/95	QC-19 x100 ERA 419	989 136	1000 135	N/A 111 - 160	99 101
Zinc	01/19/ 9 5 01/19/ 9 5	QC-19 x100 ERA 419	1009 121	1000 113	N/A 93 – 133	101 107

2 % 344

Table 2.44 Results of the MS/MSD Analysis for Water Samples WA # 0-098 Pier Drum

METAL	CLIENT#					RED CONC	% RECO		RPD R	ECOMMENI	DED
		CONC. ug/l	ug/l	Dup. ug/l	Spike ug/l	Dup. u g/ 1	Spike	Dup.		⊔MIT % Rec	RPD
	A00636 B00635		2222 2222	2222 2222	3307 3222	3356 3184	93 96	96 95	2 2	75-125 75-125	20 20
	A00636 B00635	0.600 ND	55.6 55.6	55.6 55.6	68.8 58.3	64.4 62.4	123 105	115 112	7 7	75-125 75-125	20 20
Arsenic	A00636	3.50	55.6	55.6	73.8	69.8	127 *	119	6	75-125	20
	B00635	2.80	55.6	55.6	60.2	60.4	103	104	Ö	75-125	20
Barium	A00636	18	556	556	564	572	98	100	1	75-125	20
Barium	B00635	17	556	556	573	564	100	98	2	75-125	20
	A00636	ND		222	221	224	99	101	1	75-125	20
Beryllium	B00635	ND	222	222	225	222	101	100	1	75-125	20
	A00636 B00635	0.0 5 0 0.070	55.6 55.6	55.6 55.6	68.0 60.4	64.0 61.6	122 109	115 111	6 2	75-125 75-125	20 20
	A00636 B00635	2. 40 1.10	55.6 55.6	55.6 55.6	63.2 55.2	58.2 57.4	109 97	100 101	9 4	75-125 75-125	20 20
Cobalt	A00636	. 3	222	222	221	226	98	100	2	75~125	20
	B00635	1	222	222	228	223	102	100	2	75-125	20
Copper	A00636	ND	222	222	220	225	99	101	2	75-125	20
Copper	B00635	1	222	222	227	224	102	100	1	75-125	20
	A00636	9431	2222	2222	11320	11360	85	87	2	75-125	20
Iron	B00635	9248	2222	2222	11270	11250	91	90	1	75–125	20
	A00636 B00635	0.400 0.100	55.6 55.6	55.6 55.6	65.6 55.0	55.1 53.7	117 9 9	98 96	18 2	75-125 75-125	20 20
								_			_
Manganese Manganese		4 3	222 222	222 222	218 223	222 220	96 99	98 98	2 1	75-125 75-125	20 20
Mercury	A00636	0.1	2.00	2.00	2.10	2.00	100	95	5	75-125	20
	B00635	0.1	2.00	2.00	2.10	2.10	100	100	Ö	75-125	20
Nickel	A00636	ND		222	220	221	99	99	0	75-125	20
Nickel	B00635	ND	222	222	224	219	101	99	2	75-125	20
	A00636	0.90	55.6	55.6	55.3	51.2	98	91 97	8	75-125	20
Selenium	B00635	0.20	55.6	55.6	46.6	48.6	84	87	4	75–125	20
	A00636 B00635	0.6 0.1	55.6 55.6	55.6 55.6	47.1 47.6	61 39.4	84 86	109 71 1	26 * 19	75-125 75-125	20 20
Silver	B00033										
	A00636 B00635		55.6 55.6	55.6 55.6	37.9 35.9	38.3 38.6	68 * 65 *	69 ' 69 '		75-125 75-125	20 20
	A00636	ND		556	543	550	98	99	1	75-125	20
	B00635	ND		556	553	544	100	98	2	75-125 75-125	20
Zinc	A00636	ND	222	222	211	214	95	96	1	75-125	20
Zinc	B00635	ND	222	222	216	215	97	97	0	75-125	20

Table 2.45 Results of the Blank Spike Analysis for Water Samples WA # 0-098 Pier Drum

METAL	Spiked Conc ug/l	RECOVERED CONC ug/l	% RECOVERY	RECOMMENDED LIMIT
Aluminum	2222	2145	97	75-125
Antimony	55.6	59.8	108	75-125
Arsenic	55.6	59.1	106	75-125
		541	97	75125
Barium	556			
Beryllium	222	218	98	75–125
Cadmium	55.6	66.0	119	75-125
Calcium	2222	2284	103	75-125
Chromium	55.6	54.8	99	75-125
Cobalt	222	219	99	75–125
Copper	222	222	100	75-125
Iron	2222	2199	99	75-125
Lead	55.6	59.1	106	75-125
Magnesiun	2222	2078	94	75-125
Manganese	e 222	215	97	75-125
Mercury	2.0	2.1	105	75-125
Nickel	222	220	99	75-125
Potassium	2222	1669	75	75-125
Selenium	55.6	56.8	102	75-125
Silver	55.6	60.2	108	75-125
Sodium	2222	2115	95	75-125
Thallium	55.6	52.2	94	75-125
Vanadium	556	539	97	75-125
Zinc	222	212	95	75-125

Table 2.46 Results of the QC Standard Analysis for Soil Samples WA # 0-098 Pier Drum

METAL	DATE ANALYZED	CONTROL	CONC. RECOVERED	TRUE VALUE	95 % CONFIDEN% INTERVAL	RECOVER
	<u>-</u>	STANDARD	ug/l	ug/l		
Aluminum	01/26/95	QCS/ICP-2	1071	1000	870 - 1100	107
	01/26/95	ERA-419	229	210	150 -263	109
Antimony	01/30/95	TMMA#2	102	100	81.65-125.67	102
Arsenic	01/30/95	TMMA #1	50	50	41.9-55.9	100
Barium	01/26/95	QCS/ICP-2	1045	1000	890 - 1070	105
	01/26/95	ERA-419	204	204	167 — 241	100
Beryllium	01/26/95	QC-19 x100	1049	1000	N/A	105
	01/26/95	ERA-419	59	56	44 - 66	105
Cadmium	01/26/95	QC-19 x100	1041	1000	N/A	104
	01/26/95	ERA-419	106	98	75 – 115	108
Calcium	01/26/95	QC-19 x100	994	1000	N/A	99
Chromium	01/26/95	QC-19 x100	1076	1000	N/A	108
	01/26/95	ERA-419	1053	950	779 – 1120	111
Cobalt	01/26/95	QC-19 x100	1071	1000	N/A	107
	01/26/95	ERA-419	111	103	85 – 122	108
Copper	01/26/95	QC-19 x100	1042	1000	N/A	104
	01/26/95	ERA-419	80	77	63 ~ 91	104
ron	01/26/95	QC-19 x100	1084	1000	N/A	108
	01/26/95	ERA-419	134	125	93 – 147	107
Lead	01/26/95	QC-19 x100		1000	N/A	105
	01/26/95	ERA-419	104 *	82	65 ~ 97	127
Magnesium	01/26/95	QC-19 x100	1034	1000	N/A	103
Manganese	01/26/95	QC-19 x100	1052	1000	N/A	105
	01/26/95	ERA-419	130	122	100 - 144	107
Mercury	01/25/95	WP032 #2	4.0	3.89	2.87 - 5.02	103
Nickel	01/26/95	QC-19 x100	1078	1000	N/A	108
	01/26/95	ERA-419	103	92	75 – 110	112
Potassium	01/26/95	QCS/ICP-2	9763	10000	8750 — 11100	98
Selenium	01/31/95	TMMA #1	51	50	39.4-57.4	102
Silver	01/26/95	QCS/ICP-2	530	500	440 560	106
	01/26/95	ERA-419	42	42	29 - 49	100
Sodium	01/26/95	QCS/ICP-2	1019	1000	830 - 1470	102
Thallium	01/30/95	TMMA #2	56	50	39.9-57.97	112
√anadium		QC-19 x100		1000	N/A	103
	01/26/95	ERA-419	141	135	111 ~ 160	104
Zinc	01/26/95	QC-19 x100		1000	N/A	105
	01/26/95	ERA-419	122	113	93 ~ 133	108

Table 2.47 Results of the MS/MSD Analysis for Soil Samples WA # 0-098 Pier Drum

METAL	CLIENT#	SAMPLE (CONC. mg/kg (Spike	IAL CONC Dup. mg/kg	RECOVE Spike mg/kg	RED CONC Dup. mg/kg	% RECC Spike	DVERY Dup.	RPD
Antimony	A00643	ND	4.02	4.05	2.59	2.51	64	62	4
Arsenic	A00643	0.208	4.02	4.05	4.21	4.05	99	95	5
Barium	A00643	3.62	92.8	78.9	89.6	79.5	93	96	4
Beryllium	A00643	0.023	46.4	39.4	46.4	40.8	100	103	3
Cadmium	A00643	ND	46.4	39.4	46.8	39.9	101	101	0
Chromium	A00643	0.06	46.4	39.4	48.0	41.6	103	105	2
Cobalt	A00643	ND	46.4	39.4	47.1	40.9	101	104	2
Copper	A00643	ND	46.4	39.4	44.7	39.4	96	100	4
Lead	A00643	0.93	46.4	39.4	49.1	43.0	104	107	3
Manganese	A00643	ND	46.4	39.4	46.7	40.6	101	103	2
Mercury	A00643	ND	0.322	0.331	0.338	0.347	105	105	0
Nickel	A00643	0.16	46.4	39.4	48.1	41.9	103	106	2
Selenium	A00643	0.053	4.02	4.05	4.25	4.18	104	102	2
Silver	A00643	0.29	46.4	39.4	36.8	38.3	79	96	20
Thallium	A00643	0.010	4.02	4.05	3.88	3.98	96	98	2
Vanadium	A00643	ND	92.8	78.9	91.3	79.9	96	101	3
Zinc	A00643	ND	46.4	39.4	50.7	42.2	109	107	2

Table 2.48 Results of the Blank Spike Analysis for Soil Samples WA # 0-098 Pier Drum

METAL	Spiked Conc mg/kg	Sand Blk Conc mg/kg	RECOVERED CONC mg/kg	% RECOVERY
Aluminum	400	17.9	472	114
Antimony	5.00	ND	4.12	82
Arsenic	5.00	ND	4.85	97
Barium	100	ND	101	101
Beryllium	50.0	ND	53.1	106
Cadmium	50.0	ND	51.4	103
Calcium	400	ND	429	107
Chromium	50.0	ND	53.3	107
Cobalt	50.0	ND	53.3	107
Copper	50.0	ND	51.5	103
Iron	400	30.7	451	105
Lead	50.0	1.83	52.6	102
Magnesium	400	ND	411	103
Manganese	50.0	ND	52.4	105
Mercury	0.400	ND	0.420	105
Nickel	50.0	ND	53.8	108
Potassium	400	8.17	365	89
Selenium	5.00	ND	5.04	101
Silver	50.0	ND	50.0	100
Sodium	400	ND	405	101
Thallium	5.00	ND	4.96	99
Vanadium	100	ND	104	104
Zinc	50.0	ND	54.7	109

MORN E AFIG.

Table 2.49 Results of the QC Standard Analysis for TCLP Samples WA # 0-098 Pier Drum

METAL	DATE	CONTROL	CONC. RECOVERED	TRUE VALUE	95 % CONFIDENCE	% RECOVERY
	ANALIZED	STANDARD	ug/l	ug/l	INTERVAL	•
Arsenic	02/01/95	TMMA #1	50.8	50	41.9-55.9	102
Barium	02/01/95 02/01/95	QC-7 x100 ERA 419	990 200	1000 204	N/A 167 - 241	99 98
Cadmium	02/01/95 02/01/95	QC-19 x100 ERA-419	1014 104	1000 98	N/A 75 - 115	101 106
Chromium	02/01/95 02/01/95	QC-19 x100 ERA-419	1031 1014	1000 950	N/A 779 - 1120	103 107
Lead	02/02/95	TMMA #1	53.9	50	43.4~56.3	108
Mercury	01/20/95	TMMA #1	1.9	2	1.40-2.49	95
Selenium	02/01/95	TMMA #1	50.7	50	39.4-57.4	101
Silver	02/01/95 02/01/95	QC-7 x100 ERA 419	992 42	1000 42	N/A 29 -49	99 100

Table 2.50 Results of the MS/MSD Analysis for TCLP Samples WA # 0-098 Pier Drum

METAL	CLIENT#	SAMPLE CONC.	-	RIGINAL CC Dup.	RECO\ Spike	ERED CONC. Dup.	% RECC	VERY Dup.	RPDR	ECOMMENI LIMIT	DED
		ug/l		ug/l	ug/I	ug/l	Оржо	Dup.		% Rec	RPD
Arsenic	A00527	0.9	55.6	55.6	57.8	59.2	102	105	2	75-125	20
Barium	A00527	11	556	556	570	578	101	102	1	75-125	20
Cadmium	A00527	ND	222	222	217	219	98	99	1	75-125	20
Chromium	A00527	2	222	222	224	226	100	101	1	75-125	20
Lead	A00527	ND	55.6	55.6	35.8	36.7	64 *	66 *	2	75-125	20
Mercury	A00527	0.1	2.00	2.00	2.10	2.10	100	100	0	75-125	20
Selenium	A00527	0.1	55.6	55.6	56.2	54.3	101	98	3	75-125	20
Silver	A,D00527	3	222	222	223	226	99	100	1	75-125	20

Table 2.51 Results of the Blank Spike Analysis for TCLP Samples WA # 0-098 Pier Drum

METAL	Spiked Conc ug/l	RECOVERED CONC ug/l	% RECOVERY	RECOMMENDED LIMIT
Arsenic	55.6	59.1	106	75-125
Barium	556	551	99	75-125
Cadmium	222	215	97	75-125
Chromium	222	224	101	75-125
Lead	55.6	57.2	103	75-125
Mercury	2.00	2.10	105	75-125
Selenium	55.6	60.7	109	75-125
Silver	222	225	101	75-125

Table 2.52 Results of the QC Standard Analysis for Drum Liquid Samples WA # 0-098 Pier Drum

	DATE ANALYZED		CONC. RECOVERED ug/l		95 % CONFIDENCE INTERVAL	% RECOVERY
Juminum	02/13/95	QC-7 x100	1017	1000	N/A	102
	02/13/95	ERA 419	240	210	150 - 263	114
Intimony	02/08/95	TMMA #2	121	100	81.65-125.67	121
rsenic	02/08/95	TMMA #1	50.9	50	41.9-55.9	102
and the	02/13/95	QC-7 x100	987	1000	N/A	99
Barium	02/13/95	ERA 419	191	204	167 – 241	94
Beryllium	02/13/95	QC-19 x100	1018	1000	N/A	102
serymum.	02/13/95	ERA 419	55	56	44 – 66	98
	02/13/95	QC-19 x100	1017	1000	N/A	102
Cadmium	02/13/95	ERA 419	100	97.7	75–115	102
Calcium	02/13/95	QC-19 x100	1022	1000	N/A	102
	00/10/05	QC-19 x100	1032.0	1000	N/A	103
Chromium	02/13/95	ERA 419	979	950	779-1120	_/ 103
	00/40/05	QC-19 x100	1025	1000	N/A	103
Cobalt	02/13/95 02/13/95	ERA 419	106	103	85 - 122	103
		00 40.400	1015	1000	N/A	102
Copper	02/13/95 02/13/95	QC-19 x100 ERA 419	75	77	63 – 91	97
	02/13/95	QC-19 x100	1045	1000	N/A	105
Iron	02/13/95	ERA 419	128	125	93 – 147	102
Lead	02/14/95	TMAA#1	51	50	43.4-56.3	101
Magnesiul	m 02/13/95	QC-19 x100	1005	1000	N/A	101
_	00/40/05	QC-19 x100	1021	1000	N/A	102
Manganes	e 02/13/95 02/13/95	ERA 419	122	122	100 - 144	100
A 10 - 1 1	02/13/95	QC-19 x100	1034	1000	N/A	103
Nickel	02/13/95	ERA 419	96	92	75 – 110	104
Potassium	02/13/95	QC-7 x100	9787	10000	N/A	98
Selenium	02/08/95	TMMA #1	53	50	39.4-57.4	106
0.1	00/43/05	QC-7 x100	996.0	1000	N/A	100
Silver	02/13/95 02/13/95		40	41.8	29-49	96
Sodium	02/13/95	QC-7 x100	995	1000	N/A	100
Thallium	01/17/95	TMMA #2	52.8	50	39.9-57.97	106
	00/40/05	QC-19 x10	o 979	1000	N/A	98
Vanadiur	n 02/13/95 02/13/95		130	135		96
_	00110101	- 00 10-10	00 1012	1000	N/A	101
Zinc	02/13/95 02/13/95		120	113		106

Table 2.53 Results of the Blank Spike Analysis for Drum Liquid Samples WA # 0~098 Pier Drum

METAL	Spiked Conc ug/l	RECOVERED CONC ug/l	% RECOVERY	RECOMMENDED LIMIT
Aluminum	2000	1971	99	75-125
Antimony	50.0	57.0	114	75-125
Arsenic	50.0	54.9	110	75-125
Barium	500	485	97	75-125
Beryllium	200	199	100	75-125
Cadmium	200	191	96	75-125
Calcium	2000	3069	153	75-125
Chromium	200	203	102	75-125
Cobalt	200	196	98	75-125
Copper	200	202	101	75-125
Iron	2000	2028	101	75-125
Lead	50.0	58.7	117	75-125
Magnesium	2000	1986	99	75-125
Manganese	200	198	99	75-125
Nickel	200	204	102	75-125
Potassium	2000	2107	105	75-125
Selenium	50.0	56.6	113	75-125
Silver	200	190	95	75-125
Sodium	2000	2033	102	75-125
Thallium	50.0	53.2	106	75-125
Vanadium	500	490	98	75-125
Zinc	200	491 **	246 *	75-125

^{**} High recovery could be contaminated

Table 2.54 Results of the Blank Spike Analysis for the Drum Solid Samples WA # 0~098 Pier Drum

2 1. 34.13

	Spike	Sample	BS		QC Limits
Parameter	Added	Conc.	Recov.	%	
	(ug/L)	(ug/L)	(ug/L)	Recovery	Rec.
Aluminum	10000	ND	9640	96	75-125
Antimony	50	ND	48.5	97	75-125
Arsenic	25	ND	23.8	95	75-125
Barium	5000	ND	5180	104	75-125
Beryllium	4.0	ND	3.53	88	75-125
Cadmium	10.0	ND	9.6	96	75-125
Calcium	1000	ND	950	95	75-125
Chromium	1000	ND	810	81	75-125
Cobalt	1000	ND	1030	103	75-125
Copper	1000	ND	980	96	75-125
Iron	1000	ND	1050	105	75-125
Lead	25	ND	25.5	102	75-125
Magnesium	250	ND	250	100	75-125
Manganese	1000	ND	1110	111	75-125
Mercury	5.0	ND	4.82	96	75-125
Nickel	1000	ND	1040	104	75-125
Potassium	1000	ND	1050	105	75-125 🗸
Selenium	25	ND	25.1	100	75-125
Silver	1000	ND	970	97	75-125
Sodium	250	ND	260	104	75-125
Thallium	25	ND	22.6	90	75-125
Vanadium	10000	ND	10300	103	75-125
Zinc	250	ND	260	104	75-125

Table 2.55 Results of the MS/MSD Analysis for the TCLP Analysis of the Drum Solid Samples WA # 0 \sim 098 Pier Drum

Sample ID: 00911						2	6	345
	Spike	Sample	MS		MSD			QC
Analyte	Added	Conc.	Recov.	%	Recov.	%		Limits
	(ug/L)	(ug/L)	(ug/L)	Recovery	(ug/L)	Recovery	RPD	RPD Rec.
Aluminum	10000	ND	8280	83	8440	84	2	20 75-125
Antimony	50	ND	39.3	79	37.1	74 *	6	20 75-125
A rsen ic	25	ND	25.4	102	24.5	98	4	20 75-125
Barium	5000	ND	5370	107	5370	107	0	20 75-125
Beryllium	4.0	ND	4.72	118	4.50	113	5	20 75-125
Cadmium	10.0	167	182	NC	182	NC	NC	20 75-125
Calcium	1000	435000	447000	NC	438000	NC	NC	20 75-125
Chromium	1000	ND	910	91	900	90	1	20 75-125
Cobalt	1000	470	1290	82	1290	82	0	20 75-125
Copper	1000	ND	960	98	980	98	0	20 75-125
ron	1000	2190	3080	89	3100	91	2	20 75-125
_ead	25	102	136	NC	138	NC	NC	20 75-125
Magnesium	250	4300	4400	NC	4300	NC	NC	20 75-125
Manganese	1000	350	1400	105	1410	106	1	20 75-125
Mercury	5.0	ND	4.85	97	4.94	99	2	20 75-125
Nickel	1000	ND	1030	103	1030	103	0	20 75-125
Potassium	1000	4200	5300	110	5200	100	10	20 75-125
Selenium	25	ND	16.1	64 *	15.7	63 *	3	20 75-125
Silver	1000	ND	960	96	950	95	1	20 75-125
Sodium	250	3100000	3200000	NC	2800000	NC	NC	20 75-125
Thallium	25	ND	ND	MI	ND	Mi	Mi	20 75-125
/anadium	10000	ND	7680	77	7670	77	0	20 75-125
Zinc	250	440	620	72 *	620	72 *	0	20 75-125

CCLIN

Table 2.55 Results of the MS/MSD Analysis for the TCLP Analysis of the Drum Solid Samples WA # 0-098 Pier Drum

Sample ID: 00911

	Spike	Sample	MS		MSD			1	QC
Analyte	Added	Conc.	Recov.	%	Recov.	%		Li	imits
	(ug/L)	(ug/L)	(ug/L)	Recovery	(ug/L)	Recovery	RPD	RPD	Rec
Aluminum	10000	ND	8280	83	8440	84	2	20	75 – 125
Antimony	50	ND	39.3	79	37.1	74 *	6	20	75 - 125
Arsenic	25	ND	25.4	102	24.5	98	4	20	75 - 125
Barium	5000	ND	5370	107	5370	107	0	20	75 - 125
Beryllium	4.0	ND	4.72	118	4.50	113	5	20	75 - 125
Cadmium	10.0	167	182	NC	182	NC	NC	20	75-125
Calcium	1000	435000	447000	NC	438000	NC	NC	20	75-125
Chromium	1000	ND	910	91	900	90	1	20	75-125
Cobalt	1000	470	1290	82	1290	82	0	20	75-125
Copper	1000	ND	960	98	980	98	0	20	75-125
ron	1000	2190	3080	89	3100	91	2	20	75-125
_ead	25	102	136	NC	138	NC	NC	20	75-125
/lagnesium	250	4300	4400	NC	4300	NC	NC	20	75-125
Manganese	1000	350	1400	105	1410	106	1	20	75-125
Mercury	5.0	ND	4.85	97	4.94	99	2	20	75-125
lickel [*]	1000	ND	1030	103	1030	103	0	20	75-125
otassium	1000	4200	5300	110	5200	100	10	20	75 - 125
Selenium	25	ND	16.1	64 *	15.7	63 *	3	20	75 - 125
Silver	1000	ND	960	96	950	95	1	20	75-125
Sodium	250	3100000	3200000	NC	2800000	NC	NC	20	75-125
hallium	25	ND	ND	MI	ND	MI	MI	20	75-125
/anadium	10000	ND	7680	77	7670	77	0	20	75-125
linc	250	440	620	72 *	620	72 *	0	20	75 - 125

QA/QC FOR BTU, ASH, MOISTURE, SULFUR and CHLORINE

Sample 00913 was chosen for the duplicate analysis for the drum liquid samples that were analyzed for BTU, and ash. The relative percent differences (RPDs), listed in Table 2.56, were 0 (zero) and 12. Both results were within the acceptable QC limits.

Sample 00912 was chosen for the duplicate analysis for the drum liquid samples that were analyzed for moisture. The RPD, also listed in Table 2.56, was 16. The result was within the acceptable QC limit.

Non-REAC sample AB42470 was chosen for the drum liquid samples that were analyzed for sulfur. The RPD, also listed in Table 2.56, was not calculated because sulfur was not detected in either analysis.

Non-REAC sample AB 42469 was chosen for the drum liquid samples that were analyzed for chlorine. The RPD, also listed in Table 2.56, was 4. The result was within the acceptable QC limit.

Table 2.56 Results of the Duplicate Analysis for the Drum Liquid Samples WA # 0-098 Pier Drum

Parameter	Sample	Units	First Analysis	Second Analysis	RPD	QC Limit
BTU	00913	BTU/lb	5470	6177	12	20
Ash	00913	%	0.26	0.26	0	20
Water	00912	%	53	62	16	20
Sulfur	AB42470	%	ND	ND	NC	20
Chlorine	AB 42469	%	55.6	53.3	4	20

EPA Contract 68-C4-0022 REAC, Ec. Jn, NJ (908) 321-4200

CHAIL JE CUSTODY RECORD

Project Name:_

00065 Ö

SHEET NO. 6F /

Phone (20%) 3214 20/4

FOR SUBCONTRACTING USE ONLY Helels 1 4/08 **Analyses Requested** PCB/Lot \$ Container/Preservative # of Bottles Special Instructions Date Collected 10 lm 95 101298 10 1-45 ĕĕ Sample Identification Matrix DC . - S - - A Potable Water Groundwater Surface Water Sludge Sampling Location PENCH 4 RK#1 PW. GW. SW. Sample No. DE0917 00913 71 POOL Sediment Drum Solids Drum Liquids Other 01 3 OFF REAC #

FROM CHAIN OF **CUSTODY#**

Items/Reason	Relinquished By	Date	Received By	Date	Time	Items/Reason	Relinquished By	Date	Beceived By	Date	Time
3 / Andrea /	Minul Uch 3	201.95	B Jewan	156/02/1	1.50	1/ Analyss	Steva	3/18/1	(Miss routh	130/95 1437	12/2/
			B. Luin	12/05//	ر جو	1/184/18.56	8 Leven	36/8/	JAM.	14	9:30 1
		,	School	18/1/2//	6.50	1111101015	8 Sam	8/2/8	11/15 18/11	LAK STOR FOR	200
	$\gamma^{\prime\prime}(j)$	5/11/2		/ /				, ,	s. //.		
	7								1		
FORM #4											8/94

			•	S								ζ.				- 		
25. 1. 1. 1.	Mable		\\										NLY		Time	3.50		16/3
07.00	┣ ╌ ╿ ╶┤		11	\coprod			\prod	-					USEC		Date	sspet)		
5-01-55066 No: -05066 SHEET NO. LOF L	Patring		7 7		$\setminus \mid \mid$								FOR SUBCONTRACTING USE ONLY		*			T
5-6 1-8	40	- -	-		+	+	H	-	+	+	$+\!$		TRAC	PO	scatufed by	1		
	RWA		7/7		Λ								BCON	FROM CHAIN OF CUSTODY #	1/4		\coprod	
1/ 474		+	+	H	$+$ \	+	H	+	+	+	+		R SUI	FROM CHAI	Oste	1/05/1		
1 122	407		7/7										Fo	<u></u>	ed By	ġ,		
		-	+	H - H		+	H	+		+	+	1		Yes No	Relinquished By	1/2		
Program Phoras	ContainerPreservative					\mathbb{N}							eived	109	2	7	3	4
USTO MAN	in the state of th	777				'							All Samples Received	Jc Cool	ms/Resson	BNA		
OF CO	1		\perp	-			\mathbb{H}						Sampl	ا خ الا	d pema	160	19	
H. C. S. S. S. S. S. S. S. S. S. S. S. S. S.	of Bottles	444	11				1					Special Instructions:	₹		Time	057	П	$\prod_{i=1}^{n}$
			10 h			+	H	\forall				pecial ins	-	Temp. Sample	Dete	13000	\prod	
Project Name:_ Project Number RFW Contact:_ on	Date Collected	12 de 15	Sem 15									1	į				$\dagger \dagger$	+
Pro Pro RF	Date D	999	0.0					\perp				-		₹	Received By	42		
entific	Matrix	886	22						\prod				òģò	A STA	2	12		
Pro Pro RF Sample Identification	cation	ens	* 5 km									Ordehle Weist	Groundwater Surface Water		Date	270	\prod	T
	Sampling Location	333	Y									A de de	Groun	Studge Studge As the State of t	9,	7.1	$\dagger \dagger$	1
1-0022	Sam	225	NE	4				-	ig			}		्ष न्या १	Reijnquished By	111		
, NJ 00 68-C4	Sample No.	917	915 1916							\mathbb{N}			5		2	1		
REAC, Luison, NJ (908) 321-4200 EPA Contract 68-C4-0022 0 30 %	Sam	689	1600 0094				\prod	-	-				Sediment Drum Solids Drum Linuida	out of the	2 8	==		2
REAC, L. (908) 321- EPA Cont	PEAC.		33				The second of th					ll _ä		*	Terra/Reseon	14		
	\	4 27 20	BB	_111		. Seemen	ac.	a.	<u>.</u>	: 14			నే దో దే	ix '	<u>L</u>	121	لــــــــــــــــــــــــــــــــــــــ	
	,									J	ملدل							

14:23 FOR SUBCONTRACTING USE ONLY 03663 E E SHEET NO._OF__ 16.AS Date Pettho ö TAC Analyses Requested FROM CHAIN OF CUSTODY # Date Notals Relinquished By CHAIL, JE CUSTODY RECORD Phone:_ 334704000100980 Container/Preservative ر د د separate anotherio Items/Reason SONT. FNA to /motel # of Bottles Special Instructions Time 4/4 M Project Number: 1/2/K Project Name: RFW Contact: Ş. Date Collected Soil Water Oil Air Sample Identification Received By Matrix 3 Potable Water Groundwater Surface Water Sludge Sampling Location Book Date/ MW - A ハーンソイ MY -27/7 t,e14 Relinguished By EPA Contract 68-C4-0022 PW -GW -SW -SL -GHT DOBY BC00645 Sample No. EF00641 A Pot 64 4900D) REAC, Eu., Jn, NJ Sediment Drum Solids Drum Liquids Other (908) 321-4200 at Mark 1116 Items/Reason 01138 REAC # FORM #4 08. DL. SD-

00199

É Ż 1432 TITIO FOR SUBCONTRACTING USE ONLY 03870 VOA S SHEET NO._OF__ Date PCB/PesT Received By Y Ö Analyses Requested FROM CHAIN OF CUSTODY # BNA X Date Relinquished By Soil record insurance buffles water to cold on 155119. CHAIL JE CUSTODY RECORD 0334704020100160 Phone 14103 Container/Preservative 1402 14°C tems/Reason (BNA & JA 2 rada **AOB** 1/20 Time # of Bottles alci 113/KS Date 6 A- metale filtered Squirte anotymic Project Number._ RFW Contact:_ Project Name: Date Collected Muller Received By Sample Identification Matrix Potable Water Groundwater Surface Water Sludge Cate Cate 11/23) Sampling Location Relinquished By 12.30 SW -SW -SL -SHE DOGUODE EPA Contract 68-C4-0022 EF 00 640 Ch 9000 ADDO640 Foo (31 4605659 15000 24 120002 Sediment Drum Solids Drum Liquids Other 71 3063 Sample No. 12000 D 4500635 REAC, Ear Ju, NJ Items/Reason (908) 321-4200 Me. IX FAC. 2016 ٠ ×

TIMe FOR SUBCONTRACTING USE ONLY 03671 SHEET NO._OF_ 12 Act Date Ast/1855 Ö i'x, Analyses Requested FROM CHAIN OF CUSTODY # Date Stale Contraction Relinquished By JE CUSTODY RECORD 240 000 000 Phone: Container/Preservative CHAIL II MINING ttems/Reason (mater) #/V0> A- metals y. B- metals y. # of Bottles Separte analysis Special Instructions: Time JUHU Project Number: 3/6/1/19/0 Project Name: RFW Contact: Date Date Collected Received By Sample Identification Michael Very Matrix , , , o Ł Sampling Location Potable Water Groundwater Surface Water Sludge Date $\lambda_{p, +}$ Relinquished By EPA Contract 68-C4-0022 PW -GW -SW -SL -HI COUSE ABNO 636 OFF ONLOSE 4500073 70000 D00(158 EFOCUS8 Sample No. 4600638 REAC, Eu n, NJ (908) 321-4200 Sediment Drum Solids Drum Liquids Other Items/Reason REAC # 260 97 X 55-B F Matrix: ۰. × 00.307

Project Name: Project Number RFW Contact: Sample Identification	Sampling Location Matrix Date Collected # Of Bottos	2	PW - Potable Water S - Soil Special Instructions: Sw - Groundwater W - Water FA - Alf R - Alf	CONDUCATION SAND 18 - MADE - AND AND STAND STAND STANDS INCOMESSARE (STAND) SAND STANDS STAND
ec. Jn, NJ 1-4200 ntract 68-C4-0022	REAC# Sample No. Sampling Location REAC# Sample No. Sampling Location Sample No. Sampling Location Sample No. Sampling Location Sample No. Sampling Location Sample No. Sampling Location Sample No. Sampling Location Sample No. Sample No. Sampling Location Sample No. Sample No. Sampling Location Sample No. Sample No. Sampling Location Sample No. Sample No. Sampling Location Sample No. Sample No. Sampling Location Sample No. Sample No. Sampling Location Sample No. Sample No. Sampling Location Sample No. Sampling Location Sample No. Sample No. Sampling Location Sample No. Sample No. Sample No. Sampling Location Sample No. S	1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	Itrix: Sediment BW- Drum Solids GW- Drum Liquids SW- SL-	iquots removed to any control of some sound to the real of the rea

6 3.20 1500 FOR SUBCONTRACTING USE ONLY Time SHEET NO. 24011 No: 10027 Date になると WEST Best **ANALYSES REQUESTED** FROM CHAIN OF PAceived By CUSTODY # **CORD/LAB WORK REQUEST** BNA BNA Date Relinquished By Phone - ¥4243 11- Poly/42 HNG L Poly/HND3 L Amber/42 Poly/HN03 Ambar/4°c LAMPO/4c Amball'c 03347-040-0078-01 6/Pethea. But Container/ Preservative Items/Reason J/natals Soil Special Instructions: water 00527 D, E, F = M3/MSD Matrix Date Collected Bottles প্ত Time CHAIN OF CUSTODY RFW Contact By Hold or MESS 1/18/15 Date Project Number SAMPLE IDENTIFICATION Project Name:--Sampling Location 13/45/ A00642 Date A00643 44900 Surface Water Sludge Potable Water Groundwater Roy F. V. ston, Inc. REAC, Edison, N.J. EPA Contract 68-03-3482 Ξ Relinquished By D. Holdana PW . GW . SW . CO0526 A,000527 B,E00527 C00525 A00526 F 0052 A00525 B00525 B00526 Sample No. Andysis | Drum Liquids COS - Drum Solids Items/Reason . Other REAC # L. IIIX W



Roy F. Weston, Inc. GSA Raritan Depot Building 209 Annex (Bay F) 2890 Woodbridge Avenue Edison, New Jersey 08837-3679 908-321-4200 • Fax 908-494-4021

January 17,1995

Attn:

Joe Larusso

Project # 3347-040-001-0098, Pier Drum

As per Weston REAC Purchase Order number 08-31480, dated 01/13/95, please analyze samples according to the following parameters:

Analysis/Method	Matrix	# of samples					
VOA/CLP compounds	Soil	3					
TCLP VOA/1311	Soil	3					
VOA/CLP compounds	Drum solids	5					
BNA/CLP compounds	Drum solids	5					
Pest/PCB/CLP compounds	Drum solids	5					
Metals/CLP compounds	Drum solids	5					
Corrosivity/SW-846 Chapter 7	Drum solids	5					
Ignitability/SW-846 Chapter 7	Drum solids	5					
Reactivity/SW-846 Chapter 7	Drum solids	5					
Ash Content/ASTM D-189	Drum solids	5					
% Sulfur/ASTM D-129-64	Drum solids	5					
% Chlorine/ASTM D-808-81	Drum solids	5					
Total Organic Halogens/SW-846 9020 A	Drum solids	5					
BTU/ASTM D-240-76/2015-77	Drum solids	5					
TCLP VOA, BNA, Pest/PCB, Metals/1311 Analyze for Compounds in CLP.	Drum solids	5					
Data package including diskette deliverables as per attached Deliverables Requirements							

2 ... 747

Samples are expected to arrive at your laboratory on January 13, 1994. All applicable QA/QC (MS/MSD) analysis will be performed on each of our sample matrix. Preliminary results tables including MS/MSD's plus a signed copy of our Chain of Custody is due at REAC 10 business days from sample receipt, with the complete data package due 21 business days from time of sample receipt. The complete data package must include all items on the deliverables checklist. If your laboratory cannot meet the delivery date, please give best delivery date possible.

ALL ORGANIC EXTRACTIONS ON SOLIDS IE: BNA, PEST/PCBMUST BE BY SOXHLET EXTRACTION

Should any questions or problems arise concerning this project, please call John Johnson at (908) 321-4248 or fax to (908) 494-4020. Any contractual question, please call Cindy Snyder at (908) 321-4296. Thank you

Sincerel

George Armstrong

Data Validation and Report Writing Group Leader

Roy F. Weston, Inc. / REAC Project

GA: jj Attachments

CC.

R. Singhvi

G. Powell

Central File

0098\non\mem\9501\sub\0098con

V. Kansal

Subcontracting File

M. Van Clef

C. Snyder

R. Lewis

G. Armstrong